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Second Quarter 2013 Groundwater Monitoring Report

**Former Powerine Refinery
12345 Lakeland Road, Santa Fe Springs, CA**

**SLIC No. 0318, ID No. 2040071
CAO 97-118**

Prepared on Behalf of

**Isola Law Group, LLP
Lodi, California**

Prepared for

**Regional Water Quality Control Board
Los Angeles Region**

Prepared By

The logo for MUREX environmental, inc. features a stylized red bull head icon to the left of the word "MUREX" in a bold, serif font. Below "MUREX" is the word "environmental, inc" in a smaller, lowercase sans-serif font. The entire logo is set against a white background.

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1.0 INTRODUCTION

On behalf of Isola Law Group, LLP, Murex Environmental (Murex) has prepared this *Second Quarter 2013 Groundwater Monitoring Report* for the former Powerine Refinery property located at 12345 Lakeland Road in Santa Fe Springs, California (Site; **Figure 1**).

1.1 Purpose

The objective of the quarterly groundwater monitoring is to evaluate groundwater quality beneath the site and adjacent properties (**Figure 2**) and to provide regular updates to the Regional Water Quality Control Board, Los Angeles Region (RWQCB). This report presents the groundwater monitoring activities performed between March 29, 2013 and April 15, 2013, in accordance with the RWQCB Cleanup and Abatement Order (CAO) No. 97-118.

1.2 Site Description and History

The Site is approximately 55 acres in size and is bordered to the north by Florence Avenue, to the south by Lakeland Road, and to the east by Bloomfield Avenue (**Figure 2**). Commercial/light industrial properties border the site to the west. The site was operated as an oil refinery from the 1930s until July 1995. Historical aerial photographs indicate that the western portion of the site may have been used for agricultural purposes from approximately 1928 to 1938. Oil production-related structures such as ponds and aboveground holding tanks may have also been located onsite during this time period (Haley & Aldrich, Inc. [Haley & Aldrich], 2005). The refinery is not currently in operation; however, some of the refinery structures remain onsite. These structures are scheduled to be removed prior to the redevelopment of the property for commercial/light industrial use.

Previous refining operations included processing crude oil into several grades of fuel including kerosene, leaded gasoline and aviation fuel, unleaded gasoline, jet fuel, high and low-sulfur diesel, fuel oil, and petroleum coke. Soil and groundwater quality beneath and in proximity to the site have been impacted by past site operations. Soil and groundwater investigations are being conducted pursuant to a CAOs (No. 97-118) issued by the RWQCB to Powerine Oil Company (CENCO Refining Company) in 1997 (Haley & Aldrich, 2005).

2.0 GROUNDWATER SAMPLING ACTIVITIES

Quarterly groundwater monitoring has been conducted since August 1986. The previous monitoring event was performed by Murex in January and February 2013. The following subsections summarize work completed during the second quarter 2013 monitoring event.

2.1 Monitoring Network

The quarterly groundwater monitoring program currently includes the existing 59 wells, as listed in **Table I** and shown on **Figure 2**. These wells include:

- Twenty-two on-Site groundwater monitoring wells: MW-101, MW-103, MW-104A, MW-105, MW-201, MW-202, MW-204, MW-205, MW-504, MW-701, MW-702, MW-703, MW-704, MW-705, MW-706, W-9, W-10, W-11, W-12, W-17A, W-17B, and W-17C;
- Twenty-five down-gradient off-Site groundwater monitoring wells of which:
 - Four are located on the former Lakeland (aka "Coaster") property: MW-501A, MW-502, MW-503B, and MW-707; and
 - Twenty-one are located on the Metropolitan State Hospital (MSH) property: MW-600A, MW-601A, MW-603, MW-604, MW-605, MW-606, MW-607, MW-708, MW-709, MW-710, MW-711, MW-712, MW-713, MW-714, MW-715, W-14A, W-14B, W-14C, W-15A, W-15B, and W-15C;
- Seven off-Site groundwater monitoring wells located to the southeast on the Walker property including: EW-1, W-1, W-3A, W-4, W-16A, W-16B, and W-16C;
- Three off-Site groundwater monitoring wells located to the east on the Bloomfield property that include: MW-106A, MW-107A, and MW-203; and
- Two on-Site, deep, former water production wells identified as W-7 and W-8.

2.2 Groundwater Gauging

Murex inspected and measured the depth to groundwater in all 59 of the wells on March 29, 2013. During gauging, wells are also checked for the presence and thickness of free-phase petroleum hydrocarbons (FPPH) product. Of those, 19 wells were dry, and three wells contained free-phase petroleum hydrocarbon (FPPH).

Table II summarizes the groundwater elevation and free product thickness measurements.

2.3 Free-Phase Petroleum Hydrocarbon (FPPH) Measurements

Wells that initially exhibit the presence of FPPH are purged until they become dry or until approximately 6 to 10 well volumes are evacuated. Thereafter, the wells are inspected for the return of FPPH and if present, its thickness is measured. For wells in which FPPH does not return, groundwater is sampled for analysis. Further discussion of the wells exhibiting free product is presented in **Section 3.2**.

2.4 Groundwater Purging

The groundwater monitoring wells that contained groundwater, with the exception of production wells W-7 and W-8, were purged via a dedicated vacuum stinger that was connected to a truck-mounted vacuum pump truck operated by Nieto & Sons. W-7 and W-8 are deep production wells and are sampled without purging water from them first. During purging, extracted groundwater volume and quality were recorded. The parameters measured during purging were temperature, pH, electrical conductivity, dissolved oxygen (DO), oxidation-reduction potential (ORP), color, and odor. The results of the field parameter testing are summarized in **Table IV**. Purged groundwater was disposed of by Nieto & Sons at the wastewater treatment system in operation at the Site.

2.5 Groundwater Sampling and Analysis

Following purging, groundwater samples were collected by disposable bailer from the wells, placed in sample containers, stored in pre-cooled ice chests, and transported under proper chain-of-custody procedures to Sunstar Laboratories, Inc. (Sunstar Labs) of Lake Forest, California, California Department of Public Health Environmental Laboratory Accreditation Program (ELAP) #2250.

The chemicals of concern (COCs) impacting groundwater in the area of study include total petroleum hydrocarbons as gasoline (TPHg), BTEX compounds (i.e., benzene, toluene, ethylbenzene, and xylenes), and petroleum product additives (i.e., oxygenates) and breakdown byproducts, such as methyl tert-butyl ether (MTBE) and tert-butyl alcohol (TBA). Therefore, collected groundwater samples were analyzed for the following:

- TPHg by U.S. Environmental Protection Agency (USEPA) Method 8015M, and
- Volatile organic compounds (VOCs) with oxygenates by USEPA Method 8260B.

Results of these analyses are summarized in **Table III**. Results of the field-measured parameters are shown in **Table IV**.

2.6 Quality Assurance/Quality Control

In accordance with the Quality Assurance/Quality Control (QA/QC) plan, Murex collected and submitted field duplicate samples and trip blanks for laboratory analysis as a quality assurance/quality control measure.

2.6.1 Trip Blanks

Trip blanks (provided by SunStar Lab) accompanied each daily groundwater sample shipment to evaluate the potential contamination of field samples during storage and transport. Trip blanks were analyzed for VOCs only.

2.6.2 Duplicates

Duplicate samples, which assess the precision of the laboratory analyses, were collected from wells MW-503B, MW-706, MW-708, and MW-713. This represents a duplicate frequency equal to approximately 13% relative to the total number of wells sampled. The duplicates followed the same analytical protocols as their respective primary samples. The results of the duplicate analyses are included with the original sample results in **Table III**.

2.6.3 Equipment Blanks

Equipment blanks were not collected because dedicated stingers were used to purge the wells and new disposable bailers were used for sampling, therefore eliminating cross-contamination between wells during the purging and sampling process.

2.6.4 Laboratory QA/QC Program

Laboratory QA/QC reports were reviewed to confirm proper completion of data validation tests, including batch QC results, method blanks, laboratory control samples, matrix spikes, and duplicates. The results of lab QC tests were within acceptable limits.

3.0 RESULTS & DISCUSSION

This section presents the results of the second quarter 2013 groundwater monitoring event. As mentioned earlier in the report, well completion details are provided in **Table I**. Groundwater level measurements and groundwater elevations are summarized in **Table II**. Comprehensive analytical results, including historical and recent results, are compiled in **Tables III**. **Table IV** contains a summary of bio-attenuation and field-measured parameter readings.

Figure 3 shows the groundwater elevation measured at each monitoring well, as well as the overall gradient and direction of groundwater flow on-Site. **Figure 4** depicts the same information for the entire monitoring well network. **Figure 5** shows the concentrations and estimated contour lines of TPHg measured in each well, and **Figure 7** shows similar concentrations and contour lines for benzene and MTBE.

Well measurement and groundwater sampling forms are attached as **Appendix A**. Laboratory reports and completed chain-of-custody forms are included in **Appendix B**.

The presentation of the chemical testing results in this report does not distinguish between constituents in groundwater that likely originated from the Site and those that are resultant from other sources located off-Site. Chemicals in groundwater related to off-Site sources are further discussed in Section 4.3.

3.1 Groundwater Surface Elevations and Gradient

Groundwater surface elevations were calculated for each well by subtracting the water level measurement from the top of casing elevation (**Tables I and II**). Groundwater elevations were adjusted for wells containing FPPH, assumed to have a relative density of 0.80, which is typical for mean density of various petroleum hydrocarbon mixtures. Groundwater elevations, contour lines, flow direction and gradient are shown on **Figure 4**.

Based on groundwater level measurements obtained on March 29, 2013, first-encountered groundwater beneath the site vicinity ranges in elevation from 15.74 to 51.29 feet above mean sea level (ft-amsl). Wells W-7 and W-8 are former production wells, with screens situated deeper than 500 feet below ground surface (ft-bgs). Their elevations were higher, between 48.87 and 63.91, respectively.

In general, groundwater elevations were lower to those measured in the first quarter 2013 monitoring event. For the wells that are constructed to a depth of about 125 ft-bgs or less,

groundwater elevations had exhibited steady decreases for several years until the third quarter 2010, when they experienced a significant increase. By the fourth quarter 2012, the groundwater elevations apparently leveled off, and it appears that groundwater elevations in the Site vicinity are now experiencing a decreasing trend. The groundwater elevations in the wells screened deeper (greater than 125 ft-bgs and up to 200 ft-bgs) appear to indicate similar patterns to the shallower screened wells. As a whole, the average change in groundwater elevation over all the wells measured was a decrease of approximately 3.55 feet from the first quarter 2013 sampling event. **Appendix C** includes hydrographs depicting the change in groundwater elevation over time for all the wells.

The average horizontal groundwater gradient is approximately 0.008 foot per foot (ft/ft), as shown in **Figure 4**, which was similar to the previous monitoring period, and represents what is considered a moderately steep gradient. The groundwater flow direction originates from the northeast and turns south across the area of study. This flow direction is relatively consistent with those historically reported in previous investigations.

3.2 Free-Phase Petroleum Hydrocarbons

Measurable FPPH, also known as light non-aqueous-phase liquid or LNAPL, was detected in monitoring wells EW-1, W-15A, and MW-708 (**Table II**). Well W-15A continues to exhibit FPPH, which was first measured in 2011. FPPH was measured at a thickness of 0.88 feet in EW-1, 2.51 feet in W-15A, and 0.27 feet in MW-708. During previous monitoring events going back many years, FPPH was also historically detected in wells MW-101, MW-103, MW-104, MW-201, MW-202, MW-203, MW-204, MW-205, MW-206, MW-501, MW-502, MW-503, MW-503B, MW-504, MW-600, MW-600A, MW-601, MW-601A, W-3A. The majority of these wells are now dry.

Over the previous two years, groundwater samples were collected from wells containing FPPH, after the FPPH was removed through vacuum purging. This exercise produced qualitative data for wells that would not have been sampled otherwise. The data resulting from these samples is considered imprecise given the possibility that minute quantities of FPPH were present in the sample collected. Having collected several rounds of sample data in 2011 and 2012, Murex plans to return to the previous operating procedure whereby wells containing FPPH are not sampled.

3.3 Groundwater Analysis

Within the area of study, COCs impacting groundwater include TPHg, BTEX, MTBE, and TBA. The following text presents the sampling results of the current groundwater monitoring event. Laboratory analytical results are summarized in **Table III**; laboratory reports and

completed chain-of-custody forms are included in **Appendix B**. Field-measured parameters are presented in **Table IV**.

3.3.1 Chemicals of Concern (COCs)

Gasoline is the major release product associated with the Site and is present in the Site's groundwater monitoring network as FPPH. Constituents of gasoline include BTEX compounds, in addition to oxygenated additives and breakdown byproducts, such as MTBE and TBA. The analytical result of each COC was compared to the established California Maximum Contaminant Level (MCL) in drinking water, as applicable. The following table presents the MCL for each COC, as well as the minimum and maximum concentrations detected for the current monitoring event.

Constituent	MCL (µg/L)	Q2 2013 Monitoring Event	
		Minimum Value (µg/L)	Maximum Value (µg/L)
TPHg	N/A ¹	96	25,000
Benzene	1	0.7	5,000
Toluene	150	0.61	1,500
Ethylbenzene	300	0.58	510
Xylenes (total)	1,750	0.56	2,720
MTBE	13	2.1	800
TBA	12 ²	19	350

¹ Not applicable – MCL not established for TPHg

² California Notification Level (former Action Level) – MCL not established for TBA

The analytical results for the current monitoring event indicate that concentrations of COCs within the area of study are generally similar to previous monitoring events. Elevated concentrations of TPHg, BTEX, MTBE, naphthalene, and 1,2,4-trimethylbenzene are apparent in the southwest portion of the Site, and in a dissolved-phase plume that extends south to the central portion of the MSH property. In addition to the aforementioned constituents, this plume also contains elevated concentrations of TBA and 1,3,5-trimethylbenzene. To the east (cross- to up-gradient) of the Site's monitoring network, elevated concentrations of TPHg, benzene, naphthalene, and TBA are present, though the source of these constituents is unknown.

Analytical results and isoconcentration contours are presented in **Figure 5** (TPHg), **Figure 6** (benzene), and **Figure 7** (MTBE).

3.3.2 Other VOCs

In addition to the aforementioned COCs, several additional VOCs were detected in groundwater during this monitoring event. Some of these compounds, such as naphthalene, n-propylbenzene, and trimethylbenzene, for instance, are related to petroleum hydrocarbon releases.

Conversely, also detected were chlorinated solvents, such as tetrachloroethylene (PCE), trichloroethene (TCE), 1,1-dichloroethane (1,1-DCA), 1,1-dichloroethene (1,1-DCE), 1,2-dichloroethane (1,2-DCA), and cis-1,2-dichloroethene (cis-1,2-DCE), among others, which we believe are the result of off-Site contamination entering the Powerine well network.

The U.S. EPA and the RWQCB are aware of the chlorinated solvents in groundwater through their oversight of the cleanup of a Superfund site located to the north, and up-gradient of the Site. Murex provides this data to the U.S. EPA on a periodic basis.

3.3.3 Biodegradation Parameters

Biodegradation of TPHg most commonly occurs by aerobic, nitrate-reducing, ferric iron (Fe^{3+})-reducing, sulfate-reducing, or methanogenic respiration. TPHg and BTEX serve as electron donors for microbial metabolism in aerobic biodegradation. Electron acceptors include oxygen, nitrate, Fe^{3+} , sulfate, and carbon dioxide.

In general, if sufficient oxygen is present, aerobic biodegradation will occur first. When DO concentrations fall below approximately 0.5 mg/L (an anoxic environment), denitrification will begin if nitrate is present. After most nitrate has been consumed, Fe^{3+} reduction will begin if Fe^{3+} is present. Fe^{3+} concentrations will decrease, while Fe^{2+} concentrations will increase. After most Fe^{3+} is consumed, sulfate reduction will begin if sulfate is available. After most sulfate has been consumed, methanogenesis, which involves carbon dioxide as an electron acceptor, begins. During methanogenesis, methane concentrations increase (Department of the Navy, 1998).

The results discussed below indicate that biodegradation, whether aerobic or anaerobic, may be occurring in the local environment around the wells that were sampled for biodegradation parameters.

3.3.3.1 Field-Measured Parameters

Field pH, DO, and oxidation-reduction potential (ORP) data were collected from 33 monitoring wells using an YSI 556 water quality meter (**Table IV**). The meter was inserted into grab water samples, collected from the vacuum truck intake during well purging.

Wells MW-104A, MW-106A, and W-10 did not generate enough groundwater volume during purging sufficient enough to collect water quality parameters during sampling.

- **pH** – This parameter quantifies the acidity or alkalinity of a solution. Results ranged from 7.55 to 8.45, indicating a neutral to slightly alkaline environment that is suitable for the growth of alkalophilic bacteria and microorganisms that thrive at a circumneutral pH.

- **DO** – Oxygen is the preferred electron acceptor in the biodegradation of petroleum hydrocarbons. When aerobic biodegradation occurs, DO concentrations are expected to decline as microorganisms use the electron acceptor during respiration. The vacuum stinger method used to purge the wells introduces oxygen into the groundwater. Therefore, DO data is not representative of the actual oxygen content. It is likely very low in wells exhibiting higher TPH concentrations, since oxygen is the first compound used up in the biological degradation of petroleum.

- **ORP** – This parameter is a measure of electron activity, which reflects the oxidizing or reducing nature of the environment. ORP values are generally negative under reducing conditions (gaining electrons) and positive under oxidizing conditions (losing electrons). Negative ORP values were observed in 27 of the 33 wells measured.

ORP values ranged from -251.8 mV in well MW-107A to 88.1 mV in Well MW-709. **Figure 8** presents concentrations for ORP.

Hydrogen sulfide (produced from the reduction of sulfate in groundwater, after oxygen is used up) was detected during purging of wells exhibiting elevated TPH concentrations and low or negative ORP values, which is consistent with our understanding of the conceptual site model, and indicates that aerobic degradation of the hydrocarbons has stalled due to dissolved oxygen limitations. It is likely that the introduction of air (via bioventing for example) will enhance the process of stimulating the aerobic degradation of the constituents of concern at the site.

3.3.4 QA/QC

Duplicate sample results are provided with their primary sample results in **Tables III**. The results show similar concentrations of the analytes of interest as in their respective primary samples, as would be expected for an ELAP-certified laboratory, with the following exception.

A duplicate sample was collected at well MW-708. It should be noted that the results of TPHg for the two samples were identified as exhibiting a greater than 10% variance (25 mg/L and 240 mg/L). The concentration of 240 mg/L is not considered representative due to its abnormally elevated result. As discussed in Section 3.2, the presence of FPPH in well MW-708 prior to purging is likely the cause of this variance, since a small amount of FPPH could have entered the well between the time that the two samples were collected. The concentration of TPHg in well MW-708 is therefore simply considered to be elevated, likely between the 25 mg/L and 240 mg/L results.

Trip blank samples did not indicate the presence of VOCs, which indicates proper sample storage and confirms a lack of cross-contamination during transport, with the following exception. The trip blank submitted with samples collected on April 15, 2013 exhibited a concentration of naphthalene at 1.2 µg/L. However, the samples that accompanied this trip blank, MW-704 and MW-706, did not show elevated concentrations of naphthalene compared to recent/historical results for these wells. Naphthalene was detected at a concentration of 57 µg/L in the sample collected from well MW-704, which is similar to the concentration detected in the previous monitoring event (44 µg/L), and is in line with historical results for this well that range from 3.6 µg/L to 150 µg/L over the last several monitoring events. Naphthalene was not detected at or above its laboratory reporting limit of 1.0 µg/L in neither the initial sample collected from well MW-706 nor its duplicate sample. It should also be noted that the trip blank sample was prepared and provided by the analytical laboratory. Therefore, the detection of naphthalene in the trip blank sample is considered a laboratory cross-contaminant and did not appear to affect the results of the accompanying field samples.

Laboratory method blanks did not indicate the presence of VOCs, which indicates that laboratory detection equipment did not exhibit cross-contamination.

Laboratory control and laboratory spike samples exhibited results within acceptable limits, indicating no matrix interference and that the detection equipment was working properly.

4.0 SUMMARY & CONCLUSIONS

Groundwater monitoring was performed at and in the vicinity of the former Powerine refinery in April 2013 as part of an ongoing groundwater monitoring plan intended to evaluate chemical impacts, contaminant sources, and overall groundwater quality. This groundwater monitoring event included inspecting/gauging water levels in 59 wells and collecting samples from 40 of those wells for analysis of TPHg and VOCs.

4.1 Groundwater Surface Elevations and Gradient

A horizontal groundwater gradient of approximately 0.008 ft/ft was calculated for the second quarter 2013 groundwater monitoring event. This is consistent with historical gradient data for the site vicinity. Averaging all the wells exhibiting measurable groundwater, elevations have decreased by approximately 3.55 feet since the previous quarter. Groundwater flows from the northeast and turns due south across the area of study, which is consistent with historical measurements. Deep-screened production wells W-7 and W-8 exhibited decreases of approximately 0.5 to 1.3 vertical feet in groundwater elevation this quarter; this is likely due to the cessation of municipal water pumping operations in near proximity of the site.

4.2 Free-Phase Petroleum Hydrocarbons

Measureable free product was identified in wells EW-1, W-15A, and MW-708. These wells have exhibited FPPH in the past; although it first appeared in W-15A in 2011. The FPPH thickness measured in these wells (0.88, 2.51, and 0.27 feet, respectively) does not necessarily reflect FPPH actual thickness in the surrounding aquifer as fluctuations in water levels and permeability factors can influence FPPH accumulation in monitoring wells.

Murex has conducted a study to compare the characteristics (i.e., “fingerprints”) of FPPH samples taken from several of the monitoring wells, including wells that do not currently contain FPPH. Samples of FPPH were collected from wells W-11, MW-503B, MW-708, EW-1, and W-15A. All the samples were then submitted for fingerprinting analysis to Zymax Forensics Laboratory in Escondido, California on September 21, 2011. The findings of this study were submitted to the RWQCB on January 25, 2012 as an addendum to the June 30, 2011 FPPH Investigation Report and indicate the presence of multiple possible sources of petroleum releases that have impacted groundwater within the Powerine study area. For instance, FPPH is present in well EW-1, but TPHg and benzene concentrations are comparatively low, indicating a non-gasoline release, such as fuel oil or other petroleum product.

4.3 Groundwater Quality

The highest concentrations of TPHg detected during this sampling event were beneath the northern and central portions of the MSH (**Figure 5**). The maximum verifiable detected concentrations of TPHg were 25 mg/L in well MW-711, 8.3 mg/L in well MW-713, and 5.8 mg/L in well MW-712 (while the sample collected from MW-708 exhibited a TPHg concentration of 25 mg/L, it also contained FPPH prior to purging; therefore, the quantification of TPHg in the dissolved phase is considered potentially erroneous). Wells MW-708, MW-711, MW-712, and MW-713 are located south of the Coaster property in the north-central portion of the MSH.

Benzene, toluene, ethylbenzene, xylene, and other compounds associated with petroleum hydrocarbons largely mimic TPHg in their presence and relative concentrations in the areas associated with the plume. The maximum concentration of benzene was detected in well MW-713 at 5,000 µg/L, located south of the Coaster property in the central portion of the MSH property (**Figure 6**). The maximum concentration of MTBE was also detected in well MW-713 at 800 µg/L, located in the central portion of the MSH property (**Figure 7**) at a distance of approximately 2,000 feet from the site. Several factors, including the results of the September 2011 FPPH forensics study, indicate that more than one type of hydrocarbon release is present in well W-15A.

Changes in the petroleum hydrocarbon plume may be reflective of fluctuations in groundwater elevation. Free-phase hydrocarbons, less dense than water, often remain above the water table in a "smear zone" as groundwater elevations fall. Residual impacts in the smear zone are expected to continue to cause variation in dissolved-phase concentrations and effect when and where FPPH are measured. Murex will continue to monitor the hydrocarbon plume within the well network and provided regular updates to the RWQCB through the monitoring and reporting program. Having performed an extended study of FPPH in 2011 and 2012, Murex plans to discontinue the sampling of wells that contain FPPH in future events, which is consistent with historical practice.

4.3.1 Off-Site Sources of Petroleum Hydrocarbons

In addition to historic releases from the Site, data collected from the monitoring well network (**Figures 4, 5, and 6**) exhibits evidence of other sources. Some observations that would support the presence of alternative sources are: (1) the comparatively clean appearance of FPPH in well W-15A versus the weathered or cloudy appearance of FPPH in wells EW-1, MW-503B, and MW-708; (2) the historical presence of FPPH in wells EW-1 and W-3A, which are located east and cross-gradient of the former refinery.

In order to complete characterization of the plume, Murex will be recommending the installation of additional groundwater monitoring wells in the Site's area of study, as further discussed in **Section 4.5**.

4.3.2 Discussion of Solvent Detections

Data collected from the monitoring well network (**Table III**) exhibits the presence of substances not linked to historic releases at the Site, including chlorinated solvents. The following observations were made regarding the additional detected chemicals in groundwater within the former Powerine refinery monitoring well network.

During this sampling event, elevated PCE and TCE concentrations (i.e., between 75 and 89 µg/L) were measured in well MW-710. This is consistent with previously measured high values from MW-710. Historically, these compounds were also detected in wells MW-107A, MW-701, and MW-14C.

Cis-1,2-DCE was found in 16 of the wells sampled at concentrations generally consistent with historical levels. 1,1-DCE was detected at an elevated concentration of 41 µg/L in well MW-710. Historically, wells W-14A, W-14B, and W-14C also exhibited elevated concentrations of these constituents.

The U.S. EPA and the RWQCB are aware of the chlorinated solvents in groundwater through their oversight of the cleanup of a Superfund site located to the north, and up-gradient of the Site. Murex provides this data to the U.S. EPA on a periodic basis.

4.3.3 Assessment of Vapor Risk from Groundwater Plume

At the direction of the DTSC, Murex has conducted an off-site soil gas sampling study. The results, presented to the RWQCB and DTSC in the November 7, 2011 *Off-Site Soil Gas Survey Report*, indicate that the petroleum hydrocarbon plume does not pose a threat to off-site receptors as a result of volatilization from groundwater.

4.4 Biodegradation

Intrinsic biodegradation continues to be viable, in at least some areas of the site and vicinity, based on nitrate, sulfate, Fe^{2+} , methane, alkalinity, and ORP results from previous sampling events conducted at the site. Oxygen has been depleted, as evident by the presence of hydrogen sulfide in the deep subsurface (sulfate reduction reactions result in the formation of hydrogen sulfide). Since the main limiting factor for biodegradation of petroleum hydrocarbons is oxygen, the mechanical introduction of oxygen could stimulate aerobic biodegradation of the VOCs present in groundwater.

Murex conducted pilot testing at the site to determine the appropriate remedial technology which will effectively enhance biodegradation of the constituents of concern and reduce the extent of groundwater contamination. Based on the results and data collected during pilot testing, it appears that a combination of remedial technologies would be suited for the site. The results and conclusions of this study were submitted to the RWQCB in the Pilot Testing Report dated November 21, 2011.

4.5 Additional Characterization

Murex has studied groundwater characteristics in the Site vicinity and will be recommending to the RWQCB the installation of additional groundwater monitoring wells in the area of study. The objective of this proposed work is to better define the lateral extent of free product and dissolved-phase petroleum impacts in the vicinity of the Site. In addition, Murex intends to identify where and to what extent petroleum impacts may be resultant from sources other than the Site, and where multiple plumes, if applicable, have become comingled.

5.0 REFERENCES

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6.0 CLOSING

I certify under penalty of law that this document and all enclosures were prepared under my direction or supervision in accordance with a system designed to assure that qualified personnel properly gather and evaluate the information submitted. The information contained herein is, to the best of my knowledge and belief, true, accurate and complete, however, is reliant upon public agency records, which could be incomplete or inaccurate beyond our control.

Should you have any questions or concerns regarding the material herein, please do not hesitate to contact the undersigned at (714) 508-0800.

Sincerely,
MUREX ENVIRONMENTAL, INC.

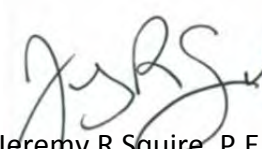

Jeremy R Squire, P.E.
Senior Engineer



Table I
Well Construction Details
Former CENCO Refinery
Santa Fe Springs, CA

Well Installation					Completion Data															Location	Reference(s) ¹	
Well ID	Date	By	Elevation		Hole Diameter (in)	Casing Diameter (in)	Screen		Depth (ft)						Elevation (ft)							
			Ground Surface (ft)	Top of Casing (ft-amsl)			Slot (in)	Length (ft)	Sand Pack		Slotted		Total Depth		Sand Pack		Slotted		Total Depth			
									Top	Bottom	Top	Bottom	Casing	Hole	Top	Bottom	Top	Bottom	Casing			Hole
Groundwater Monitoring Wells																						
EW-1	1989	Emcon	146.85	146.85	-	4	-	-	-	-	-	-	113.5	-	-	-	-	-	-	-	Walker	Versar (2000)
MW-101	8/28/1985	IT	145.19	138.00	12	4	-	20	69.5	90	70	90	90	95	66	45	65	45	45	40	Refinery	IT (1986); Versar (2000); ARCADIS (2003)
MW-103	8/30/1985	IT	137.18	139.36	12	4	-	20	-	-	79	99	99	99.5	-	-	58	38	-	37	Refinery	IT (1986); Versar (2000); ARCADIS (2003)
MW-104	8/24/1985	IT	-	-	12	4	-	20	-	-	76.5	96 5	97	99	-	-	66	46	-	43	Refinery	IT (1986); Versar (2000); ARCADIS (2003)
MW-104A	6/1999	Versar	142.38	144.13	-	4	-	-	-	-	65	100	100	-	-	-	-	-	-	-	Refinery	Versar (2000); measured well depth
MW-105	12/1995	TriHydro	-	141.16	-	4	-	-	-	-	68	98	98	100	-	-	-	-	-	39	Refinery	Versar (2000); measured well depth
MW-106	12/1995	TriHydro	-	-	-	4	-	-	-	-	74	104	106.45	106	-	-	-	-	42	42	Bloomfield	Versar (2000)
MW-106A	2/20/2006	N&M	152.92	152.81	8	4	0.02	27	82	110	83	110	110	110	70	42	69	42	42	42	Bloomfield	Well completion report
MW-107	12/1995	TriHydro	-	-	-	4	-	-	-	-	75	105	107 55	108	-	-	-	-	41	41	Bloomfield	Versar (2000)
MW-107A	2/20/2006	N&M	147.37	147.02	8	4	0.02	27	82	110	83	110	110	110	64	36	63	36	36	36	Bloomfield	Well completion report
MW-201	9/10/1985	IT	134.86	135.65	12	4	-	30	66	103	72	102	102	103	67	30	61	31	31	30	Refinery	IT (1986); Versar (2000); ARCADIS (2003)
MW-202	9/23/1985	IT	139 00*	140.62	16	4	-	30	58	105	63	93	93	105	70	23	65	35	35	23	Refinery	IT (1986); Versar (2000); ARCADIS (2003)
MW-203	9/13/1985	IT	144.08	143.71	12	4	-	30	64.7	107	77	107	107	119	78	36	66	36	36	24	Bloomfield	IT (1986); Versar (2000); ARCADIS (2003)
MW-204	9/19/1985	IT	141.15	142.90	12	4	-	30	67.5	105	73.3	103.3	103.3	105	73	35	67	37	37	35	Refinery	IT (1986); Versar (2000); ARCADIS (2003)
MW-205	9/14/1985	IT	140 00*	140.09	12	4	-	30	65.5	103	69.5	99 5	99.5	104.5	73	35	69	39	39	34	Refinery	IT (1986); Versar (2000); ARCADIS (2003)
MW-206 ²	9/18/1985	IT	-	-	-	4	-	30	62.5	104	71	101	101	104	67	26	59	29	29	26	Lakeland	IT (1986); Versar (2000); ARCADIS (2003)
MW-501	6/9/1986	IT	-	-	-	4	-	30	-	-	71	101	101	107	-	-	58	28	-	22	Lakeland	IT (1986); Versar (2000); ARCADIS (2003)
MW-501A	3/1999	ATC	131.26	130.89	-	4	-	-	-	-	75	95	95	95	-	-	-	-	-	35	Lakeland	Versar (2000); measured well depth
MW-502	6/11/1986	IT	131.88	131.00	-	4	-	30	-	-	74	104	104	104	-	-	54	24	-	24	Lakeland	IT (1986); Versar (2000); ARCADIS (2003)
MW-503	6/13/1986	IT	-	-	-	4	-	30	-	-	80.5	110.5	110.5	111	-	-	51	21	-	20	Lakeland	IT (1986); Versar (2000); ARCADIS (2003)
MW-503B	1/1999	Versar	133.03	132.66	-	4	-	-	-	-	69	109	109	109	-	-	-	-	-	21	Lakeland	Versar (2000); measured well depth
MW-504	6/18/1986	IT	-	137.18	-	4	-	50	-	-	58	118	95.76	118	-	-	77	17	-	17	Refinery	IT (1986); Versar (2000); ARCADIS (2003)
MW-600	8/15/1990	ENSR	-	-	-	4	-	30	-	-	78	108	108	110	-	-	42	12	-	10	MSH	IT (1986); Versar (2000); ARCADIS (2003)
MW-600A	6/1999	Versar	123.28	124.26	-	4	-	-	-	-	-	-	92.7	100	-	-	-	-	-	20	MSH	Versar (2000); measured well depth
MW-601	8/17/1990	ENSR	-	-	-	4	-	30	-	-	85	115	115	117	-	-	40	10	-	8	MSH	IT (1986); Versar (2000); ARCADIS (2003)
MW-601A	6/1999	Versar	-	-	-	4	-	-	-	-	65	100	100	100	-	-	-	-	-	27	MSH	Versar (2000); measured well depth
MW-603	12/1995	TriHydro	121.40	120.95	-	4	-	-	-	-	70	100	100	100	-	-	-	-	-	19	MSH	Versar (2000); measured well depth
MW-604	12/1995	TriHydro	140.52	140.07	-	4	-	-	-	-	73	103	103	103	-	-	-	-	-	35	MSH	Versar (2000); measured well depth
MW-605	12/1995	TriHydro	117.40	116.82	-	4	-	-	-	-	65	95	95	95	-	-	-	-	-	20	MSH	Versar (2000); measured well depth
MW-606	12/1995	TriHydro	116.90	116.06	-	4	-	-	-	-	70	100	100	100	-	-	-	-	-	14	MSH	Versar (2000); measured well depth
MW-607	12/1995	TriHydro	128.92	128.28	-	4	-	-	-	-	77	107	107	107	-	-	-	-	-	19	MSH	Versar (2000); measured well depth
W-1	12/1995	TRC	145.19	144.81	-	4	-	-	-	-	70	129	129	130	-	-	-	-	-	13	Walker	IT (1986); Versar (2000)
W-2 ²	12/1995	TRC	-	-	-	4	-	-	-	-	84	129	129	129	-	-	-	-	-	-	Walker	IT (1986); Versar (2000)
W-3 ²	12/1995	TRC	-	-	-	4	-	-	-	-	82	122	122	124	-	-	-	-	-	-	Walker	IT (1986); Versar (2000)
W-3A	-	-	137.18	136.79	-	4	-	-	-	-	-	-	111.52	115	-	-	-	-	-	21	Walker	Versar (2000)
W-4	12/1995	TRC	143.18	142.56	-	4	-	-	-	-	79	129	130	-	-	-	-	-	-	-	Walker	IT (1986); Versar (2000)
W-9	8/22/2006	Arcadis	140.37	139.84	8	2	0.01	35	73	111	75	110	110	120.5	66	28	64	29	29	19	Refinery	ARCADIS BBL (2006)
W-10	8/21/2006	Arcadis	141.39	140.71	8	2	0.01	35	73	111	75	110	110	130	67	29	65	30	30	10	Refinery	ARCADIS BBL (2006)
W-11	8/25/2006	Arcadis	141.96	142.10	8	2	0.01	35	73	111	75	110	110	119	68	30	66	31	31	22	Refinery	ARCADIS BBL (2006)
W-12	8/23/2006	Arcadis	142.93	145.15	8	2	0.01	35	75	114	75	114	114	120.5	69	30	69	30	30	24	Refinery	ARCADIS BBL (2006)
W-14A	1/22/2008- 1/30/2008	Arcadis	115.23	114.71	9	2	0.02	45	67	112	67	112	112	200	48	3	48	3	3	-85	MSH	ARCADIS (2008)
W-14B			115 00*	114.78	9	2	0.02	10	157	167	157	167	167	200	-42	-52	-42	-52	-52	-85		
W-14C			115 00*	114.78	9	2	0.02	10	185	195	185	195	195	200	-70	-80	-70	-80	-80	-85		

Table I
Well Construction Details
Former CENCO Refinery
Santa Fe Springs, CA

Well Installation					Completion Data															Location	Reference(s) ¹	
Well ID	Date	By	Elevation		Hole Diameter (in)	Casing Diameter (in)	Screen		Depth (ft)						Elevation (ft)							
			Ground Surface (ft)	Top of Casing (ft-amsl)			Slot (in)	Length (ft)	Sand Pack		Slotted		Total Depth		Sand Pack		Slotted		Total Depth			
									Top	Bottom	Top	Bottom	Casing	Hole	Top	Bottom	Top	Bottom	Casing			Hole
W-15A	11/27/2007-12/10/2007	Arcadis	127.91	127.59	10	2	0.02	45	78	126	80	125	125	200	50	2	48	3	3	-72	MSH	ARCADIS (2008)
W-15B			128 00*	127.61	10	2	0.02	10	143	156	145	155	155	200	-15	-28	-17	-27	-27	-72		
W-15C			128 00*	127.59	10	2	0.02	10	188	200	190	200	200	200	-60	-72	-62	-72	-72	-72		
W-16A	10/24/2007-10/30/2007	Arcadis	147.89	147.60	10	2	0.02	45	76	125	78	123	123	200	72	23	70	25	25	-52	Walker	ARCADIS (2008)
W-16B			148 00*	147.68	10	2	0.02	10	143	156	152	162	162	200	5	-8	-4	-14	-14	-52		
W-16C			148 00*	147.67	10	2	0.02	10	184	200	186	196	196	200	-36	-52	-38	-48	-48	-52		
W-17A	1/31/2008-2/8/2008	Arcadis	141.60	141.38	9	2	0.02	45	63	108	63	108	108	200	78	33	78	33	33	-59	Refinery	ARCADIS (2008)
W-17B			142 00*	141.37	9	2	0.02	10	159	169	159	169	169	200	-18	-28	-18	-28	-28	-59		
W-17C			142 00*	141.38	9	2	0.02	10	190	200	190	200	200	200	-49	-59	-49	-59	-59	-59		
MW-701	12/6/2010	Murex	136.87	139.48	12	4	0.02	50	77	130	80	130	130	130	59.87	6.87	56.87	6.87	6.87	6.87	Refinery	Murex (2011)
MW-702	12/15/2010	Murex	140.90	140.12	12	4	0.02	50	77	130	80	130	130	130	63.90	10 90	60.90	10.90	10.90	10.90	Refinery	Murex (2011)
MW-703	12/10/2010	Murex	134.73	137.23	12	4	0.02	50	77	130	80	130	130	130	57.73	4.73	54.73	4.73	4.73	4.73	Refinery	Murex (2011)
MW-704	12/14/2010	Murex	137.93	137.66	12	4	0.02	50	77	130	80	130	130	130	60.93	7.93	57.93	7.93	7.93	7.93	Refinery	Murex (2011)
MW-705	12/13/2010	Murex	139.16	141.94	12	4	0.02	50	77	130	80	130	130	130	62.16	9.16	59.16	9.16	9.16	9.16	Refinery	Murex (2011)
MW-706	12/9/2010	Murex	139.68	139.30	12	4	0.02	50	77	130	80	130	130	130	62.68	9.68	59.68	9.68	9.68	9.68	Refinery	Murex (2011)
MW-707	12/23/2010	Murex	128.86	128.43	12	4	0.02	50	77	130	80	130	130	130	51.86	-1.14	48.86	-1.14	-1.14	-1.14	Getty Drive	Murex (2011)
MW-708	1/12/2011	Murex	126.73	126.26	12	4	0.02	50	77	130	80	130	130	130	49.73	-3.27	46.73	-3.27	-3 27	-3.27	MSH	Murex (2011)
MW-709	1/26/2011	Murex	140.48	139.78	12	4	0.02	50	77	130	80	130	130	130	63.48	10.48	60.48	10.48	10.48	10.48	MSH	Murex (2011)
MW-710	1/13/2011	Murex	122.15	121.99	12	4	0.02	50	77	130	80	130	130	130	45.15	-7.85	42.15	-7.85	-7 85	-7.85	MSH	Murex (2011)
MW-711	1/17/2011	Murex	128.09	127.84	12	4	0.02	50	77	130	80	130	130	130	51.09	-1.91	48.09	-1.91	-1 91	-1.91	MSH	Murex (2011)
MW-712	1/24/2011	Murex	123.57	123.31	12	4	0.02	50	77	130	80	130	130	130	46.57	-6.43	43.57	-6.43	-6.43	-6.43	MSH	Murex (2011)
MW-713	1/19/2011	Murex	128.42	128.15	12	4	0.02	50	77	130	80	130	130	130	51.42	-1.58	48.42	-1.58	-1 58	-1.58	MSH	Murex (2011)
MW-714	1/20/2011	Murex	129.07	128.87	12	4	0.02	50	77	130	80	130	143	130	52.07	-0.93	49.07	-0.93	-13 93	-0.93	MSH	Murex (2011)
MW-715	1/27/2011	Murex	116.66	116.22	12	4	0.02	50	77	130	80	130	130	130	39.66	-13.34	36.66	-13.34	-13 34	-13.34	MSH	Murex (2011)
Former Groundwater Production Wells																						
					-	-	-	80	-	-	450	530	690	-	-	-	-	-	-	-	Refinery	IT (1986)
W-7	-	-	-	141.97	-	-	-	90	-	-	600	690	-	-	-	-	-	-	-	-	Refinery	
W-8	-	-	-	141.11	-	-	-	-	-	-	-	-	994	-	-	-	-	-	-	-	Refinery	

NOTES:

¹Sources: IT, 1986; Versar, 2000; Arcadis, 2003, 2006, 2008, and 2009; Dan Herlihy Environmental Services, 2006 (as shown).

²Well abandoned
ft Feet
in Inches
MSH Metropolitan State Hospital Property
amsl Above mean sea level
TOC Top of casing
* Value retrieved from Google Earth

Table II
Summary of Groundwater Level Measurements
Former CENCO Refinery
Santa Fe Springs, CA
2Q2013

Well ID	Date	Total Depth (ft)	Depth to Groundwater (ft)	Depth To FPPH (ft)	FPPH Thickness (ft)	Top of Casing Elevation (ft amsl)	Groundwater Elevation (ft amsl)
EW-1	1/18/2013	113.31	106.62	105.74	0.88	146.85	40.93
W-1	1/18/2013	129.58	109.21			144.81	35.60
W-3A	1/18/2013	111.73	DRY			136.79	NA
W-4	1/18/2013	129.50	110.43			142.56	32.13
W-7	1/18/2013	NM	93.10			141.97	48.87
W-8	1/18/2013	NM	77.20			141.11	63.91
W-9	1/18/2013	110.37	90.93			139.84	48.91
W-10	1/18/2013	110.18	97.98			140.71	42.73
W-11	1/18/2013	110.05	98.97			142.10	43.13
W-12	1/18/2013	116.10	103.66			145.15	41.49
W-14 A	1/18/2013	111.85	94.78			114.71	19.93
W-14 B	1/18/2013	167.00	93.69			114.78	21.09
W-14 C	1/18/2013	195.00	93.91			114.78	20.87
W-15 A	1/18/2013	125.40	113.59	111.08	2.51	127.59	16.01
W-15 B	1/18/2013	155.60	111.53			127.61	16.08
W-15 C	1/18/2013	197.34	111.85			127.59	15.74
W-16 A	1/18/2013	122.86	112.37			147.60	35.23
W-16 B	1/18/2013	160.25	117.92			147.68	29.76
W-16 C	1/18/2013	196.30	117.70			147.67	29.97
W-17 A	1/18/2013	108.07	97.25			141.38	44.13
W-17 B	1/18/2013	169.60	107.01			141.37	34.36
W-17 C	1/18/2013	200.00	107.08			141.38	34.30
MW-101	1/18/2013	90.72	DRY			138.00	NA
MW-103	1/18/2013	94.70	DRY			139.36	NA
MW-104A	1/18/2013	100.14	92.84			144.13	51.29
MW-105	1/18/2013	100.47	DRY			141.16	NA
MW-106A	1/18/2013	110.08	104.24			152.81	48.57
MW-107A	1/18/2013	109.29	104.03			147.02	42.99
MW-201	1/18/2013	101.60	DRY			135.65	NA
MW-202	1/18/2013	92.55	DRY			140.62	NA
MW-203	1/18/2013	102.30	DRY			143.71	NA
MW-204	1/18/2013	103.10	DRY			142.90	NA
MW-205	1/18/2013	98.27	DRY			140.09	NA
MW-501A	1/18/2013	93.27	DRY			130.89	NA
MW-502	1/18/2013	100.59	DRY			131.00	NA
MW-503B	1/18/2013	108.72	101.60			132.66	31.06
MW-504	1/18/2013	95.76	DRY			137.18	NA
MW-600A	1/18/2013	92.70	DRY			124.26	NA
MW-601A	1/18/2013	89.90	DRY			126.53	NA
MW-603	1/18/2013	97.60	DRY			120.95	NA
MW-604	1/18/2013	103.20	DRY			140.07	NA
MW-605	1/18/2013	93.98	DRY			116.82	NA
MW-606	1/18/2013	99.05	DRY			116.06	NA
MW-607	1/18/2013	107.05	DRY			128.28	NA
MW-701	1/18/2013	132.65	99.91			139.48	39.57
MW-702	1/18/2013	130.00	99.52			140.12	40.60
MW-703	1/18/2013	130.00	101.17			137.23	36.06
MW-704	1/18/2013	129.81	102.90			137.66	34.76
MW-705	1/18/2013	133.39	103.90			141.94	38.04
MW-706	1/18/2013	130.13	100.48			139.30	38.82
MW-707	1/18/2013	130.77	98.46			128.43	29.97
MW-708	1/18/2013	130.00	97.82	97.55	0.27	126.26	28.66
MW-709	1/18/2013	130.00	109.65			139.78	30.13

Table II
Summary of Groundwater Level Measurements
Former CENCO Refinery
Santa Fe Springs, CA
2Q2013

Well ID	Date	Total Depth	Depth to Groundwater	Depth To FPPH	FPPH Thickness	Top of Casing Elevation	Groundwater Elevation
MW-710	1/18/2013	130.00	96.25			121.99	25.74
MW-711	1/18/2013	130.00	103.00			127.84	24.84
MW-712	1/18/2013	130.00	99.95			123.31	23.36
MW-713	1/18/2013	130.00	105.59			128.15	22.56
MW-714	1/18/2013	132.50	106.18			128.87	22.69
MW-715	1/18/2013	134.00	97.98			116.22	18.24

NOTES:

ft Feet
FPPH Free-phase petroleum hydrocarbons
amsl Above mean sea level
NM Not measured, inaccessible
NA Not available/applicable

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
EW-1	ug/L	11/1/1989	9800	730	16	1400A								<5		9.8			<5	<5	29
EW-1	ug/L	3/1/1990		1800	300	1800								<25		<50			<25	<25	<100
EW-1	ug/L	4/1/1990		1300	290	1600								<1		20	110		<10	<10	<20
EW-1	ug/L	8/21/1998	5000	230	<50	630			<50		150	<50	<50	<50		<50	<50		<50	<50	<100
EW-1	ug/L	1/28/1999	7900	110	<50	540			<50		130	<50	<50	<50		<50	<50		<50	<50	<100
EW-1	ug/L	7/19/1999	8000	110	<25	1000			<25		<250	<25	25	<25		<25	<25		<25	<13	<13
EW-1	ug/L	1/13/2000	NS	NS	NS	NS			NS		NS	NS	NS	NS		NS	NS		NS	NS	NS
EW-1	ug/L	7/31/2000	NS	NS	NS	NS			NS		NS	NS	NS	NS		NS	NS		NS	NS	NS
EW-1	ug/L	2/6/2001	NS	NS	NS	NS			NS		NS	NS	NS	NS		NS	NS		NS	NS	NS
EW-1	ug/L	7/26/2001	NS	NS	NS	NS			NS		NS	NS	NS	NS		NS	NS		NS	NS	NS
EW-1	ug/L	5/6/2002	NS	NS	NS	NS			NS	NS	NS	NS	NS	NS		NS	NS		NS	NS	NS
EW-1	ug/L	9/25/2002	NS	NS	NS	NS			NS	NS	NS	NS	NS	NS		NS	NS		NS	NS	NS
EW-1	ug/L	11/10/2006	4800	65	<4	68	16	<4	<10	<100	42	6.9	<4	<4		8.4	6.3		<4	<4	<10
EW-1	ug/L	2/9/2007	4100	41	<2	39	9.4	<2	<5	<50	26	5.1	2.3	<2		7.8	6.5		<2	<2	<5
EW-1	ug/L	5/10/2007	3300	19	1.5	15	3.7	<4	<10	17	10	2.6	1.4	<4		6.9	6.9		<4	<4	<10
EW-1	ug/L	8/10/2007	3200	36	2.3	14	4.7	0.64	<5	15	20	3.2	1.4	<2		9.9	11		0.35	<2	<5
EW-1	ug/L	2/8/2008	4100	73	1.9	4.9	<4	<4	<10	31	5.3	0.48	<4	<4		14	9.8		0.54	<4	2.6
EW-1	UG/L	2/3/2011	4500	20	1.5	27	13	<0.50	<1.0	<10	42	<1.0	<1.0	<1.0	1.3	5.9	4.0	<1.0	<1.0	<0.50	<1.0
EW-1	UG/L	2/3/2011	4200	20	1.4	27	13	<0.50	<1.0	<10	22	<1.0	<1.0	<1.0	1.1	5.1	3.5	<1.0	<1.0	<0.50	<1.0
EW-1	UG/L	4/13/2011	4700	29	3.2	51	28	0.74	<1.0	<10	67	1.9	<1.0	<1.0	3.7	8.9	8.6	<1.0	<1.0	<0.50	<1.0
EW-1	UG/L	11/13/2012	2900	<0.50	<0.50	5.8	1.4	<0.50	<1.0	<10	120	1.3	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
EW-1	UG/L	1/29/2013	4500	<0.50	3.0	6.1	18	9.3	<1.0	<10	110	20	4.2	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
EW-1	UG/L	4/10/2013	1400	<0.50	<0.50	2.1	<1.0	<0.50	<1.0	<10	88	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-104A	ug/L	7/19/1999	<500	<0.5	<1	<1			<1		<10	<1	<1	<1		<1	5.6		<1	1.2	<0.5
MW-104A	ug/L	1/13/2000	<500	<0.5	<1	<1			<1		<10	<1	<1	<1		<1	6.7		<1	<0.5	5.7
MW-104A	ug/L	8/2/2000	<500	<0.5	<1	<1			<1		<10	<1	<1	<1		<1	5.4		<1	<0.5	<0.5
MW-104A	ug/L	2/7/2001	<500	<0.5	<1	<1			<1		<10	<1	<1	<1		<1	4.2		<1	<0.5	<0.5
MW-104A	ug/L	7/25/2001	<100	<0.5	<1	<1			<1		<10	<1	<1	<1		<1	3.9		<1	<0.5	<0.5
MW-104A	ug/L	5/7/2002	100	<0.5	<1	<1			<1	31000	<10	<1	<1	<1		<1	4.3		<1	<0.5	<0.5
MW-104A	ug/L	9/24/2002	<100	<0.5	<1	<1			<1	20000	<10	<1	<1	<1		1.4	5.4		<1	<0.5	<0.5
MW-104A	ug/L	6/30/2004	<200	<5	<5	<5			<5	30J		<5	<5	<5		2J	8.1		<5	<5	<5
MW-104A	ug/L	10/7/2005	<100	<0.5	<1	<1	<1	<1	<1	83	<10	<1	<1	<1		<1	3.4		<1	<0.5	<0.5
MW-104A	ug/L	2/15/2006	<50	<1	<5	<5	<5	<5	<1	30	<5	<5	<5	<5		<5	2		<5	<5	<5
MW-104A	ug/L	2/7/2007	540	<2	<2	<2	<2	<2	<5	120	<5	<2	<2	<2		<2	<2		<2	<2	<5
MW-104A	ug/L	5/8/2007	33	<2	0.37	<2	<2	<2	<5	340	<5	<2	<2	<2		<2	1.8		<2	<2	<5
MW-104A	ug/L	8/8/2007	<50	<2	<2	<2	<2	<2	<5	150	<5	<2	<2	<2		0.51	2.9		<2	<2	<5
MW-104A	ug/L	11/5/2007	<30	<0.28	<0.36	<0.25	<0.6	<0.3	<0.32	81	<0.41	<0.23	<0.26	<0.32		0.71	4		<0.27	<0.28	<0.3
MW-104A	ug/L	2/4/2008	<50	<2	<2	<2	<2	<2	<5	71	<5	<2	<2	<2		0.91	5.2		<2	<2	<5
MW-104A	ug/L	1/16/2009	46	<2	<2	<2	1	<2	<5	23	<5	0.55	<2	<2		0.57	4.6		<2	<2	<5
MW-104A	ug/L	4/22/2009	<50	<2	<2	<2	<2	<2	<5	38	<5	<2	<2	<2		0.62	4.5		<2	<2	<5
MW-104A	UG/L	3/3/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	3.7		<1.0	<0.50	<1.0
MW-104A	UG/L	8/4/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	4.5		<1.0	<0.50	<1.0
MW-104A	UG/L	11/3/2010	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	3.6	<1.0	<1.0	<0.50	<1.0
MW-104A	UG/L	2/2/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.1	<1.0	<1.0	<0.50	<1.0
MW-104A	UG/L	2/2/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.2	<1.0	<1.0	<0.50	<1.0
MW-104A	UG/L	4/14/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	1.0	<1.0	6.4	<1.0	<1.0	<0.50	<1.0
MW-104A	UG/L	8/24/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	3.3	<1.0	<1.0	<0.50	<1.0

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
MW-104A	UG/L	11/10/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.4	<1.0	<1.0	<0.50	<1.0
MW-104A	UG/L	11/10/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.6	<1.0	<1.0	<0.50	<1.0
MW-104A	UG/L	2/9/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	3.8	<1.0	<1.0	<0.50	<1.0
MW-104A	UG/L	5/9/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	18	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	4.3	<1.0	<1.0	<0.50	<1.0
MW-104A	UG/L	8/27/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	3.6	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-104A	UG/L	11/6/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.7	<1.0	<1.0	<0.50	<1.0
MW-104A	UG/L	1/28/2013	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-104A	UG/L	4/5/2013	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	1.0	<1.0	4.4	<1.0	<1.0	<0.50	<1.0
MW-106A	ug/L	8/2/2006	310	2.6	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		21	13		<2	<2	10
MW-106A	ug/L	11/9/2006	82	<2	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		17	14		<2	<2	7
MW-106A	ug/L	2/8/2007	270	2.6	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		20	15		<2	<2	13
MW-106A	ug/L	5/10/2007	210	1.5	<2	0.28	<2	<2	<5	20	<5	<2	<2	<2		12	9.9		0.6	<2	7.9
MW-106A	ug/L	8/9/2007	270	1.6	<2	0.6	<2	<2	<5	19	0.69	<2	<2	<2		14	12		0.83	<2	12
MW-106A	ug/L	11/7/2007	240	1.4	<0.36	0.84	<0.6	<0.3	<0.32	20	1.6	<0.23	<0.26	<0.32		9.5	11		0.7	<0.28	9.9
MW-106A	ug/L	2/5/2008	220	1.6	<2	0.42	<2	<2	<5	16	1.8	<2	<2	<2		7.8	10		0.73	<2	10
MW-106A	ug/L	1/19/2009	220	0.46	<2	<2	<2	<2	<5	17	<5	<2	<2	<2		11	13		0.99	<2	6.3
MW-106A	ug/L	4/23/2009	290	1.9	<2	3.7	<2	<2	<5	18	0.93	<2	<2	<2		6.3	5.5		0.82	<2	10
MW-106A	UG/L	3/5/2010	590	8.4	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		2.0	3.5		<1.0	<0.50	<1.0
MW-106A	UG/L	5/13/2010	460	8.6	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		2.0	<1.0		<1.0	<0.50	21
MW-106A	UG/L	8/6/2010	450	12	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		3.5	1.0		1.2	<0.50	25
MW-106A	UG/L	11/4/2010	630	0.64	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.5	<1.0	<1.0	<0.50	8.8
MW-106A	UG/L	2/3/2011	570	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-106A	UG/L	4/19/2011	480	0.63	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.0	<1.0	<1.0	<0.50	6.9
MW-106A	UG/L	8/25/2011	540	0.51	<0.50	<0.50	<1.0	<0.50	<1.0	26	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	4.8
MW-106A	UG/L	11/14/2011	440	0.87	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-106A	UG/L	2/3/2012	440	2.7	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	11
MW-106A	UG/L	5/8/2012	630	7.1	<0.50	0.87	1.5	<0.50	<1.0	13	7.2	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	23
MW-106A	UG/L	8/24/2012	470	4.8	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	11
MW-106A	UG/L	11/6/2012	610	6.9	<0.50	0.83	<1.0	<0.50	<1.0	<10	1.5	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	78
MW-106A	UG/L	1/28/2013	250	5.7	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-106A	UG/L	4/4/2013	480	6.9	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	14
MW-107A	ug/L	8/2/2006	770	3.7	<2	<2	3.4	<2	<5	<50	<5	<2	<2	<2		2.4	3.9		<2	<2	<5
MW-107A	ug/L	11/9/2006	780	24	<2	4.7	9.1	<2	<5	<50	<5	<2	<2	<2		5.3	6.2		<2	<2	<5
MW-107A	ug/L	2/8/2007	500	80	<2	21	25	<2	<5	<50	7.4	<2	<2	<2		7.4	9.6		<2	<2	<5
MW-107A	ug/L	5/10/2007	670	42	1	14	17	<2	<5	21	6	<2	0.29	<2		6	6.6		<2	<2	2
MW-107A	ug/L	8/9/2007	1000	61	2	15	41	<2	<5	18	8.5	<2	0.33	<2		9.5	8.8		0.31	<2	2.3
MW-107A	ug/L	11/7/2007	1500	44	4.2	16	26	<0.3	<0.32	35	11	<0.23	0.49	<0.32		9.4	6.4		0.3	<0.28	4.4
MW-107A	ug/L	2/5/2008	2800	19	3	3	12	<2	<5	37	3.9	<2	0.38	<2		9.2	5.6		0.29	<2	5
MW-107A	ug/L	1/19/2009	1100	13	1.9	1.5	9.9	0.43	<5	66	1.1	<2	0.29	<2		7.3	6.8		<2	<2	2
MW-107A	ug/L	1/19/2009	1200	12	1.9	1.6	9.6	0.38	<5	62	1.3	<2	0.27	<2		7.5	7.2		<2	<2	1.8
MW-107A	ug/L	4/23/2009	1300	74	1.1	13	94	0.47	<5	67	6.6	3.2	2.8	<2		10	8.5		<2	<2	1.3
MW-107A	ug/L	4/23/2009	2400	79	1.2	13	91	0.47	<5	66	7.5	3	2.7	<2		11	9.4		<2	<2	1.3
MW-107A	UG/L	3/5/2010	1100	17	0.68	1.6		<0.50	<1.0	<10	6.0	<1.0	<1.0	<1.0		7.6	6.8		<1.0	<0.50	<1.0
MW-107A	UG/L	3/5/2010	1300	16	0.66	1.7		<0.50	<1.0	<10	5.6	<1.0	<1.0	<1.0		7.4	6.4		<1.0	<0.50	<1.0
MW-107A	UG/L	5/13/2010	1500	7.6	11	4.1		2.0	4.7	<10	3.3	2.0	<1.0	<1.0		4.7	4.8		<1.0	<0.50	<1.0
MW-107A	UG/L	5/13/2010	1100	8.8	11	4.2		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		5.9	5.9		<1.0	<0.50	<1.0

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
MW-107A	UG/L	8/6/2010	1300	120	150	39		1.3	<1.0	<10	24	1.9	<1.0	<1.0		7.5	10		<1.0	<0.50	<1.0
MW-107A	UG/L	8/6/2010	1300	120	160	39		1.3	<1.0	<10	29	1.9	<1.0	<1.0		7.0	9.5		<1.0	<0.50	<1.0
MW-107A	UG/L	11/4/2010	1400	39	11	16	29	<0.50	<1.0	<10	4.1	<1.0	<1.0	<1.0	7.5	5.8	7.7	<1.0	<1.0	<0.50	<1.0
MW-107A	UG/L	11/4/2010	1600	36	10	14	26	<0.50	<1.0	<10	4.2	<1.0	<1.0	<1.0	7.1	5.1	6.9	<1.0	<1.0	<0.50	<1.0
MW-107A	UG/L	2/3/2011	740	4.1	2.2	3.2	14	<0.50	<1.0	<10	1.2	<1.0	<1.0	<1.0	3.3	2.4	3.2	<1.0	<1.0	<0.50	<1.0
MW-107A	UG/L	4/19/2011	1200	2.4	0.90	1.2	4.7	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	5.4	3.6	5.0	<1.0	<1.0	<0.50	<1.0
MW-107A	UG/L	4/19/2011	1200	2.6	0.99	1.2	5.2	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	5.9	4.2	5.9	<1.0	<1.0	<0.50	<1.0
MW-107A	UG/L	8/25/2011	590	0.95	<0.50	<0.50	1.8	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	2.4	1.7	3.4	<1.0	<1.0	<0.50	<1.0
MW-107A	UG/L	8/25/2011	480	0.84	<0.50	<0.50	1.4	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	1.9	1.4	3.0	<1.0	<1.0	<0.50	<1.0
MW-107A	UG/L	11/14/2011	550	1.0	<0.50	<0.50	1.6	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	2.0	<1.0	4.8	<1.0	<1.0	<0.50	<1.0
MW-107A	UG/L	1/31/2012	500	0.97	0.54	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	3.6	2.6	7.8	<1.0	<1.0	<0.50	<1.0
MW-107A	UG/L	5/8/2012	710	0.78	<0.50	<0.50	<1.0	<0.50	<1.0	<10	2.1	<1.0	<1.0	<1.0	1.7	1.6	3.4	<1.0	<1.0	<0.50	<1.0
MW-107A	UG/L	8/24/2012	720	1.0	<0.50	<0.50	<1.0	<0.50	<1.0	11	<1.0	<1.0	<1.0	<1.0	2.5	1.8	3.4	<1.0	<1.0	<0.50	<1.0
MW-107A	UG/L	11/6/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-107A	UG/L	1/28/2013	450	<0.50	<0.50	1.2	8.3	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-107A	UG/L	4/4/2013	180	<0.50	2.1	1.8	9.6	5.3	<1.0	<10	71	15	3.2	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-503B	ug/L	2/9/1999	10000	970	<50	420					<50	<50	<50	<50		150	110		<50	<50	<100
MW-503B	ug/L	7/19/1999	7800	630	<20	540			<20		<200	<20	<20	<20		250	180		<20	<10	<10
MW-503B	ug/L	1/14/2000	14000	1000	32	870			<20		<200	<20	<20	<20		200	210		<20	<10	<10
MW-503B	ug/L	8/4/2000	5600	610	19	500			<10		23	<10	<10	<10		160	140		<10	<5	<5
MW-503B	ug/L	2/6/2001	5800	250	<20	320			<20		<200	<20	<20	<20		150	84		<20	<10	<10
MW-503B	ug/L	7/25/2001	5700	280	<50	230			<50		<500	<50	<50	<50		57	<50		<50	<25	<25
MW-503B	ug/L	5/9/2002	4500	81	3.5	77			<2	<20000	26	2.5	2.2	<2		23	23		<2	<1	7.7
MW-503B	ug/L	9/26/2002	3300	36	9.6	140			<1	<10000	48	2.5	3.7	<1		16	18		<1	<0.5	10
MW-503B	ug/L	7/1/2004	5900	160	37	89	42	<0.5	<5	<100	42	3J	4J	<5			3J		<5	<5	<5
MW-503B	ug/L	10/5/2005	5400	1100	<20	73	38	<20	<20	<200	<200	<20	<20	<20		<20	<20		<20	<10	<10
MW-503B	ug/L	2/14/2006	5450	331	<50	12	<250	<250	<10	<100	<50	<50	<50	<50		<50	<50		<50	<50	<50
MW-503B	ug/L	8/4/2006	4700	31	<2	3.5	2.1	2	7.6	<50	<5	<2	<2	<2		3.1	7.2		<2	<2	5.8
MW-503B	ug/L	11/10/2006	3500	26	<4	4.7	<4	<4	<10	<100	<10	<4	<4	<4		<4	4.9		<4	<4	<10
MW-503B	ug/L	2/9/2007	1600	59	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		2.2	11		<2	<2	5.4
MW-503B	ug/L	5/11/2007	1800	60	0.58	2.1	1	<2	1.3	<50	1.5	<2	0.61	<2		2.6	17		0.63	0.47	7.4
MW-503B	ug/L	8/10/2007	1800	80	0.62	1.7	1.1	<2	<5	<50	<5	0.23	0.44	<2		2	19		0.48	0.64	7.6
MW-503B	ug/L	11/8/2007	2400	270	3.6	3.7	4.7	<1.2	2.8	<20	11	<0.92	<1	<1.3		<1.1	15		<1.1	<1.1	7
MW-503B	ug/L	2/11/2008	2700	220	3.1	3.4	3.5	<8	3.4	<200	18	<8	<8	<8		1.4	21		<8	<8	6.3
MW-503B	ug/L	1/21/2009	6200	410	14	39	28	<10	<25	<250	36	<10	<10	<10		<10	<10		<10	<10	25
MW-503B	ug/L	4/27/2009	4000	210	11	24	18	2.9	2.2	<50	29	0.53	2.9	<2		<2	4.8		<2	1.2	25
MW-503B	UG/L	3/8/2010	2800	40	1.4	1.7		<0.50	2.9	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	6.7
MW-503B	UG/L	5/17/2010	2900	91	1.0	1.2		<0.50	5.1	<10	1.4	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	1.6	5.7
MW-503B	UG/L	8/9/2010	3700	270	5.3	2.4		0.65	<1.0	<10	3.4	<1.0	1.3	<1.0		<1.0	<1.0		<1.0	3.8	5.4
MW-503B	UG/L	11/8/2010	8000	690	320	180	580	170	8.2	<10	97	370	140	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	5.9
MW-503B	UG/L	11/8/2010	12000	940	440	250	800	230	9.6	<10	250	450	170	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.7	6.1
MW-503B	UG/L	2/4/2011	57000	1400	7700	2900	15000	5900	<1.0	<10	5200	15000	4400	<1.0	<1.0	<1.0	2.7	<1.0	<1.0	4.8	<1.0
MW-503B	UG/L	4/15/2011	41000	3400	3200	1800	7200	2600	9.1	63	370	2100	640	<1.0	<1.0	<1.0	1.4	<1.0	<1.0	<0.50	8.0
MW-503B	UG/L	4/15/2011	39000	2200	2500	1400	5200	2000	9.0	64	260	1800	620	<1.0	<1.0	<1.0	1.5	<1.0	<1.0	<0.50	6.9
MW-503B	UG/L	8/29/2011	13000	590	270	440	1300	670	4.4	<10	200	470	150	<1.0	<1.0	<1.0	2.7	<1.0	<1.0	<0.50	1.1
MW-503B	UG/L	11/16/2011	6700	170	160	220	550	280	<1.0	<10	170	290	96	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-503B	UG/L	1/31/2012	5400	250	120	270	580	290	<1.0	<10	150	300	57	<1.0	<1.0	<1.0	3.3	<1.0	<1.0	2.0	<1.0

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
MW-503B	UG/L	1/31/2012	5200	280	120	300	650	330	<1.0	<10	170	340	55	<1.0	<1.0	<1.0	3.5	<1.0	<1.0	2.1	<1.0
MW-503B	UG/L	5/8/2012	11000	920	170	820	1800	250	<1.0	<10	150	770	100	<1.0	<1.0	<1.0	6.0	<1.0	<1.0	0.56	2.5
MW-503B	UG/L	8/30/2012	2000	130	19	100	190	39	3.9	<10	98	120	34	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-503B	UG/L	11/5/2012	680	120	2.1	5.4	19	4.4	1.3	12	23	24	5.5	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-503B	UG/L	1/30/2013	1100	52	18	41	130	55	<1.0	<10	140	120	35	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-503B	UG/L	4/8/2013	720	64	4.3	17	47	12	2.8	20	76	39	8.7	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-503B	UG/L	4/8/2013	600	62	4.1	16	44	11	2.7	19	75	36	8.1	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-701	UG/L	2/4/2011	190	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	4.3	1.6	9.5	1.7	<1.0	<0.50	<1.0
MW-701	UG/L	4/11/2011	230	1.1	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	14	2.3	14	3.8	1.0	<0.50	6.0
MW-701	UG/L	8/30/2011	190	2.5	<0.50	<0.50	<1.0	<0.50	<1.0	19	<1.0	<1.0	<1.0	<1.0	14	2.3	9.0	3.4	<1.0	<0.50	5.2
MW-701	UG/L	8/30/2011	290	2.7	<0.50	<0.50	<1.0	<0.50	<1.0	29	<1.0	<1.0	<1.0	<1.0	11	2.0	7.7	2.8	<1.0	<0.50	4.0
MW-701	UG/L	11/16/2011	310	2.5	0.62	1.4	3.5	1.8	<1.0	<10	7.6	3.4	<1.0	1.3	13	<1.0	9.2	4.6	<1.0	<0.50	<1.0
MW-701	UG/L	2/1/2012	300	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	8.9	3.8	14	4.3	<1.0	<0.50	<1.0
MW-701	UG/L	5/11/2012	260	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	15	3.8	14	<1.0	<1.0	<0.50	5.5
MW-701	UG/L	8/31/2012	350	0.75	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	2.7	16	2.9	18	5.3	<1.0	<0.50	3.7
MW-701	UG/L	8/31/2012	340	0.94	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	2.5	15	2.8	17	5.0	<1.0	<0.50	3.5
MW-701	UG/L	11/13/2012	300	0.95	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	5.9	3.1	18	5.1	<1.0	<0.50	31
MW-701	UG/L	2/4/2013	93	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	5.4	<1.0	22	4.0	<1.0	<0.50	4.0
MW-701	UG/L	4/10/2013	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	1.8	<1.0	7.3	<1.0	<1.0	<0.50	<1.0
MW-702	UG/L	2/4/2011	2300	91	0.74	0.92	<1.0	<0.50	<1.0	<10	5.2	<1.0	1.5	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-702	UG/L	4/12/2011	910	6.3	<0.50	<0.50	<1.0	<0.50	<1.0	32	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.0	<1.0	1.3	<0.50	1.1
MW-702	UG/L	8/30/2011	260	15	<0.50	<0.50	<1.0	<0.50	<1.0	59	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.9	<1.0	<1.0	<0.50	1.1
MW-702	UG/L	11/16/2011	1400	99	0.59	0.51	<1.0	<0.50	<1.0	<10	2.9	<1.0	1.0	<1.0	<1.0	<1.0	2.5	<1.0	1.2	<0.50	<1.0
MW-702	UG/L	2/9/2012	1400	480	1.3	0.65	<1.0	<0.50	<1.0	<10	3.4	<1.0	1.2	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-702	UG/L	2/9/2012	1500	470	1.3	0.71	<1.0	<0.50	<1.0	<10	3.3	<1.0	1.3	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-702	UG/L	5/11/2012	6000	2700	2.7	1.0	1.4	0.85	<1.0	<10	4.2	<1.0	4.4	<1.0	<1.0	<1.0	1.5	<1.0	<1.0	<0.50	<1.0
MW-702	UG/L	8/31/2012	1200	88	5.9	1.8	<1.0	0.94	<1.0	<10	<1.0	<1.0	2.0	<1.0	<1.0	<1.0	1.2	<1.0	<1.0	<0.50	<1.0
MW-702	UG/L	8/31/2012	4300	72	6.2	1.9	<1.0	0.99	<1.0	<10	<1.0	<1.0	2.1	<1.0	<1.0	<1.0	1.3	<1.0	<1.0	<0.50	<1.0
MW-702	UG/L	11/13/2012	65	17	<0.50	<0.50	<1.0	<0.50	<1.0	<10	3.5	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-702	UG/L	2/4/2013	1100	16	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	39	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-702	UG/L	4/10/2013	2300	15	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.3	<1.0	1.5	<0.50	<1.0
MW-703	UG/L	2/4/2011	1300	33	1.3	5.2	2.8	<0.50	<1.0	<10	1.6	1.8	<1.0	<1.0	2.0	<1.0	18	3.6	<1.0	<0.50	<1.0
MW-703	UG/L	4/12/2011	1100	76	1.4	7.8	4.8	<0.50	1.4	<10	<1.0	2.7	<1.0	<1.0	2.6	<1.0	10	1.7	<1.0	<0.50	<1.0
MW-703	UG/L	8/30/2011	2100	170	3.4	20	8.5	<0.50	3.3	50	<1.0	2.4	1.1	<1.0	1.1	<1.0	8.7	<1.0	<1.0	<0.50	1.3
MW-703	UG/L	11/17/2011	1700	170	3.8	25	5.6	<0.50	<1.0	<10	<1.0	2.5	1.2	<1.0	<1.0	<1.0	8.8	<1.0	<1.0	<0.50	<1.0
MW-703	UG/L	11/17/2011	1400	150	3.4	21	4.7	<0.50	<1.0	<10	<1.0	2.2	1.0	<1.0	<1.0	<1.0	9.2	<1.0	<1.0	<0.50	<1.0
MW-703	UG/L	2/14/2012	470	48	0.72	1.4	1.9	<0.50	<1.0	<10	1.1	<1.0	<1.0	<1.0	2.6	1.0	28	3.0	<1.0	<0.50	2.5
MW-703	UG/L	5/11/2012	500	10	<0.50	0.55	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	14	<1.0	<1.0	<0.50	1.1
MW-703	UG/L	8/31/2012	490	39	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	12	1.5	<1.0	<0.50	1.2
MW-703	UG/L	8/31/2012	430	40	<0.50	0.52	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	13	1.5	<1.0	<0.50	1.1
MW-703	UG/L	11/14/2012	280	4.1	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	14	2.5	<1.0	<0.50	9.5
MW-703	UG/L	2/4/2013	180	13	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	2.5	<1.0	<1.0	<1.0	14	1.7	<1.0	<0.50	<1.0
MW-703	UG/L	4/10/2013	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	5.7	<1.0	<1.0	<0.50	<1.0
MW-704	UG/L	2/9/2011	27000	1800	2000	610	3600	680	210	<10	120	1200	520	<1.0	2.3	<1.0	2.5	<1.0	1.2	38	<1.0

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
MW-704	UG/L	2/9/2011	26000	1900	2400	620	3700	720	430	<10	96	1300	550	<1.0	<1.0	<1.0	2.5	<1.0	1.3	40	<1.0
MW-704	UG/L	4/13/2011	5400	170	110	200	190	68	73	<10	38	<1.0	<1.0	<1.0	<1.0	<1.0	5.6	<1.0	6.0	7.0	2.0
MW-704	UG/L	8/31/2011	11000	570	600	300	540	180	180	160	58	410	170	<1.0	<1.0	<1.0	3.8	<1.0	3.5	25	1.5
MW-704	UG/L	11/17/2011	10000	550	430	420	520	180	190	<10	37	490	210	<1.0	<1.0	<1.0	3.4	<1.0	3.9	18	<1.0
MW-704	UG/L	2/14/2012	7700	310	89	390	530	95	100	73	50	500	210	<1.0	<1.0	<1.0	5.3	<1.0	5.7	5.9	3.1
MW-704	UG/L	2/14/2012	7800	320	89	410	560	96	130	80	53	510	220	<1.0	<1.0	<1.0	4.5	<1.0	4.9	6.2	2 3
MW-704	UG/L	5/14/2012	11000	450	250	360	520	99	130	45	61	410	150	<1.0	<1.0	<1.0	2.8	<1.0	3.3	12	1 2
MW-704	UG/L	5/14/2012	9000	460	260	360	530	98	140	56	77	420	150	<1.0	<1.0	<1.0	3.0	<1.0	3.4	12	1 2
MW-704	UG/L	9/4/2012	7800	580	30	550	760	33	44	24	3.6	670	260	<1.0	<1.0	<1.0	2.4	<1.0	2.6	3.4	<1.0
MW-704	UG/L	11/14/2012	8700	2200	150	1200	1700	170	610	60	150	1000	430	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	27	2 2
MW-704	UG/L	11/14/2012	14000	1800	120	1200	1500	150	260	43	100	1100	440	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	18	2.4
MW-704	UG/L	2/5/2013	1500	390	440	73	340	110	61	<10	44	91	24	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	37	<1.0
MW-704	UG/L	4/15/2013	3900	420	29	200	300	10	97	<10	57	530	170	<1.0	<1.0	<1.0	3.0	<1.0	3.2	6.1	1.1
MW-705	UG/L	2/4/2011	3100	450	3.5	5.1	6.4	0.54	90	94	6.7	<1.0	1.3	<1.0	<1.0	<1.0	2.0	<1.0	<1.0	<0.50	<1.0
MW-705	UG/L	4/12/2011	930	55	0.87	1.7	1.6	<0.50	22	31	<1.0	1.3	<1.0	<1.0	<1.0	<1.0	3.8	<1.0	<1.0	<0.50	<1.0
MW-705	UG/L	8/31/2011	1300	79	1.4	3.3	2.3	<0.50	13	66	<1.0	1.9	1.3	<1.0	<1.0	<1.0	4.2	<1.0	<1.0	0.56	1.2
MW-705	UG/L	11/17/2011	1100	56	7.6	24	29	6.3	73	<10	38	31	9.8	<1.0	<1.0	<1.0	2.1	<1.0	<1.0	<0.50	<1.0
MW-705	UG/L	2/14/2012	410	52	1.2	7.0	7.8	0.66	250	240	3.3	8.1	3.8	<1.0	<1.0	<1.0	8.9	1.3	<1.0	<0.50	1.8
MW-705	UG/L	2/14/2012	440	49	0.86	5.6	5.7	<0.50	250	230	<1.0	5.0	2.6	<1.0	<1.0	<1.0	8.3	1.3	<1.0	<0.50	1.5
MW-705	UG/L	5/14/2012	600	27	1.2	2.8	5.6	0.76	64	49	12	5.9	2.0	<1.0	<1.0	<1.0	7.4	1.4	<1.0	<0.50	<1.0
MW-705	UG/L	5/14/2012	610	36	<0.50	2.1	5.6	<0.50	60	33	<1.0	1.1	<1.0	<1.0	1.0	<1.0	8.3	1.8	<1.0	<0.50	<1.0
MW-705	UG/L	9/4/2012	100	0.79	<0.50	<0.50	<1.0	<0.50	12	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	13	2.0	<1.0	0.51	<1.0
MW-705	UG/L	11/14/2012	100	5.1	0.56	7.9	9.9	0.94	2.1	47	22	9.7	3.2	<1.0	<1.0	<1.0	9.2	2.3	<1.0	<0.50	3.6
MW-705	UG/L	11/14/2012	100	<0.50	<0.50	<0.50	<1.0	<0.50	1.7	24	<1.0	<1.0	<1.0	<1.0	1.1	<1.0	11	2.2	<1.0	0.56	<1.0
MW-705	UG/L	2/5/2013	<50	<0.50	0.74	<0.50	1.9	0.96	<1.0	37	3.2	1.6	<1.0	<1.0	<1.0	<1.0	1.2	<1.0	<1.0	<0.50	<1.0
MW-705	UG/L	4/10/2013	140	0.97	<0.50	<0.50	<1.0	<0.50	10	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	7.8	1.5	<1.0	<0.50	2.0
MW-706	UG/L	2/4/2011	390	4.9	0.57	<0.50	<1.0	<0.50	4.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	3.6	<1.0	<1.0	<0.50	<1.0
MW-706	UG/L	4/11/2011	540	9.0	<0.50	<0.50	<1.0	<0.50	5.9	89	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	6.0	<1.0	<1.0	<0.50	2.6
MW-706	UG/L	8/31/2011	1100	25	0.86	0.65	1.9	<0.50	5.4	54	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	4.3	<1.0	<1.0	<0.50	1.9
MW-706	UG/L	11/18/2011	490	9.5	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	3.3	<1.0	<1.0	<0.50	<1.0
MW-706	UG/L	2/14/2012	350	16	<0.50	<0.50	<1.0	<0.50	4.4	16	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	4.5	<1.0	<1.0	<0.50	2.5
MW-706	UG/L	5/14/2012	1300	22	1.0	0.95	2.6	0.50	6.8	16	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	3.7	<1.0	<1.0	<0.50	1.5
MW-706	UG/L	5/14/2012	1500	23	1.0	1.0	2.6	0.53	7.0	17	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	4.0	<1.0	<1.0	<0.50	1.6
MW-706	UG/L	9/4/2012	410	12	<0.50	<0.50	1.2	<0.50	5.8	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	4.8	<1.0	<1.0	<0.50	1.2
MW-706	UG/L	11/15/2012	<50	2.6	<0.50	3.0	4.1	<0.50	6.6	110	6.1	3.0	1.2	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-706	UG/L	11/15/2012	<50	3.1	<0.50	0.86	1.1	<0.50	5.6	110	2.9	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-706	UG/L	2/5/2013	<50	<0.50	0.80	0.53	2.4	1.2	<1.0	31	4.6	2.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-706	UG/L	4/15/2013	260	5.9	<0.50	<0.50	<1.0	<0.50	2.8	54	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	4.0	<1.0	<1.0	<0.50	1.0
MW-706	UG/L	4/15/2013	250	5.1	<0.50	<0.50	<1.0	<0.50	3.2	61	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	4.9	<1.0	<1.0	<0.50	1.5
MW-707	UG/L	2/4/2011	2000	520	120	7.6	120	150	15	<10	<1.0	10	7.8	4.1	8.7	<1.0	7.0	6.9	<1.0	2.7	<1.0
MW-707	UG/L	4/8/2011	7000	1000	560	180	670	310	15	<10	26	74	27	<1.0	3.2	<1.0	8.7	1.6	<1.0	4.0	<1.0
MW-707	UG/L	9/1/2011	2200	1200	95	92	1500	170	17	46	87	160	35	<1.0	<1.0	<1.0	6.6	<1.0	<1.0	<0.50	4 6
MW-707	UG/L	11/18/2011	8300	930	120	55	1900	120	<1.0	<10	150	250	53	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-707	UG/L	2/1/2012	10000	1200	150	100	1100	96	<1.0	<10	110	220	69	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-707	UG/L	5/15/2012	9700	1000	200	82	870	74	15	12	120	190	42	<1.0	<1.0	<1.0	3.2	<1.0	<1.0	<0.50	2 3

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
MW-707	UG/L	9/4/2012	6700	1400	41	26	220	29	9.7	<10	5.2	55	26	<1.0	<1.0	<1.0	3.8	<1.0	<1.0	1.3	1.5
MW-707	UG/L	11/15/2012	310	180	11	6.6	29	9.5	2.3	<10	21	11	2.7	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-707	UG/L	2/5/2013	92	49	5.4	2.5	19	5.3	1.4	<10	27	5.4	2.2	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-707	UG/L	4/8/2013	240	92	5.6	5.2	27	5.0	2.1	<10	29	6.0	2.7	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-708	UG/L	2/4/2011	530000	1400	420	3000	8100	13	330	<10	370	2200	92	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-708	UG/L	9/1/2011	38000	1900	230	1200	2200	54	2300	2500	150	440	430	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-708	UG/L	11/18/2011	18000	1100	62	630	860	30	1000	<100	180	940	390	<10	<10	<10	<10	<10	<10	<5.0	<10
MW-708	UG/L	2/10/2012	18000	1700	74	770	1000	38	830	<10	170	1100	410	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-708	UG/L	5/15/2012	57000	870	39	550	750	18	450	120	110	430	380	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	0.86	<1.0
MW-708	UG/L	9/5/2012	17000	1400	75	710	1000	32	390	<10	160	1400	520	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-708	UG/L	11/16/2012	1000	73	0.57	5.4	9.5	0.58	3.8	55	4.0	37	13	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-708	UG/L	2/11/2013	3200	46	2.8	19	39	4.0	52	<10	79	200	62	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-708	UG/L	4/11/2013	25000	1100	54	510	920	27	790	350	290	1700	670	<1.0	<1.0	<1.0	2.4	<1.0	<1.0	<0.50	<1.0
MW-708	UG/L	4/11/2013	240000	990	54	430	890	24	670	260	680	2000	780	<1.0	<1.0	<1.0	2.4	<1.0	<1.0	<0.50	<1.0
MW-709	UG/L	2/4/2011	500	16	1.0	<0.50	4.8	1.1	2.8	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-709	UG/L	4/6/2011	580	26	0.86	0.89	4.1	0.72	4.6	<10	2.7	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-709	UG/L	9/1/2011	9900	1.1	<0.50	0.91	4.6	1.2	7.6	60	<1.0	2.4	1.2	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-709	UG/L	11/21/2011	1100	<0.50	<0.50	0.77	2.1	0.75	6.4	<10	4.6	1.4	1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-709	UG/L	2/10/2012	760	<0.50	<0.50	<0.50	<1.0	<0.50	4.4	180	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-709	UG/L	5/16/2012	920	<0.50	<0.50	<0.50	<1.0	<0.50	4.7	20	1.1	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-709	UG/L	9/5/2012	670	<0.50	0.86	<0.50	1.8	0.67	2.2	23	12	1.2	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-709	UG/L	11/16/2012	650	1.7	<0.50	<0.50	<1.0	<0.50	2.4	100	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-709	UG/L	2/11/2013	310	<0.50	<0.50	<0.50	<1.0	<0.50	3.8	86	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-709	UG/L	4/11/2013	850	<0.50	<0.50	<0.50	<1.0	<0.50	2.1	160	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-710	UG/L	2/8/2011	93	0.84	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	55	93	2.9	14	41	3.1	0.81	1.3
MW-710	UG/L	2/8/2011	110	0.75	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	54	89	2.9	14	41	3.1	<0.50	1.2
MW-710	UG/L	4/7/2011	<50	0.81	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	76	72	4.1	19	56	4.9	1.5	2.0
MW-710	UG/L	4/7/2011	100	0.84	<0.50	<0.50	<1.0	<0.50	<1.0	<10	1.0	<1.0	<1.0	82	92	4.0	18	54	4.7	1.5	1.9
MW-710	UG/L	9/2/2011	380	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	76	97	2.0	17	50	4.3	1.2	1.1
MW-710	UG/L	9/2/2011	100	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	58	<1.0	<1.0	<1.0	76	100	2.2	18	54	4.6	1.2	1.3
MW-710	UG/L	11/21/2011	95	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	51	71	1.5	13	35	3.6	<0.50	<1.0
MW-710	UG/L	11/21/2011	79	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	52	71	1.5	13	34	3.4	<0.50	<1.0
MW-710	UG/L	2/1/2012	170	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	66	110	2.1	23	71	6.0	<0.50	<1.0
MW-710	UG/L	5/16/2012	130	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	53	77	1.2	19	48	4.4	<0.50	<1.0
MW-710	UG/L	9/5/2012	100	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	3.8	1.0	<1.0	77	91	<1.0	16	56	3.9	<0.50	1.2
MW-710	UG/L	11/16/2012	95	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	81	130	2.0	19	86	4.8	<0.50	8.2
MW-710	UG/L	2/11/2013	55	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	75	100	<1.0	18	52	4.0	<0.50	<1.0
MW-710	UG/L	2/11/2013	64	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	86	110	<1.0	19	59	4.4	<0.50	1.1
MW-710	UG/L	4/12/2013	130	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	75	89	1.7	16	41	3.6	<0.50	<1.0
MW-711	UG/L	2/8/2011	11000	520	440	120	380	250	11	<10	260	180	110	<1.0	8.4	<1.0	4.5	<1.0	<1.0	<0.50	7.5
MW-711	UG/L	4/6/2011	7100	<0.50	<0.50	65	160	50	20	<10	420	52	36	<1.0	1.1	<1.0	2.6	<1.0	<1.0	<0.50	8.7
MW-711	UG/L	9/2/2011	44000	1600	1800	650	3000	1100	25	<10	620	1800	550	<1.0	<1.0	1 3	3.8	<1.0	<1.0	<0.50	17
MW-711	UG/L	11/21/2011	14000	370	290	530	1800	790	<1.0	<10	880	480	98	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-711	UG/L	2/10/2012	23000	1900	2100	440	1800	770	14	<10	360	480	150	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
MW-711	UG/L	5/16/2012	25000	2900	3200	730	3000	1200	14	<10	370	<1.0	300	<1.0	<1.0	<1.0	3.0	<1.0	<1.0	<0.50	5 9
MW-711	UG/L	9/5/2012	28000	2100	2000	640	2000	1100	5.9	<10	370	720	120	<1.0	<1.0	<1.0	2.2	<1.0	<1.0	<0.50	5 8
MW-711	UG/L	11/16/2012	35000	6200	7000	1400	4500	2300	4.4	41	350	430	210	<1.0	<1.0	5 9	19	<1.0	1.2	<0.50	120
MW-711	UG/L	2/11/2013	410	75	35	9.8	44	20	<1.0	<10	220	27	7.6	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-711	UG/L	2/11/2013	410	71	33	9.6	43	20	<1.0	<10	240	24	7.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-711	UG/L	4/12/2013	25000	2000	1500	450	2000	720	<1.0	<10	440	640	200	<1.0	<1.0	<1.0	2.5	<1.0	<1.0	<0.50	5 9
MW-712	UG/L	2/9/2011	14000	1200	520	380	1800	390	23	<10	98	460	170	<1.0	<1.0	<1.0	2.6	<1.0	<1.0	<0.50	<1.0
MW-712	UG/L	4/7/2011	94	860	140	270	1100	170	32	<10	140	580	220	<1.0	1.8	<1.0	3.4	<1.0	<1.0	0.64	2 2
MW-712	UG/L	9/2/2011	6300	440	77	100	350	72	19	<10	43	180	76	<1.0	<1.0	<1.0	2.8	<1.0	<1.0	0.71	<1.0
MW-712	UG/L	11/21/2011	8000	600	60	90	310	60	<1.0	<10	65	140	72	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-712	UG/L	2/13/2012	8300	850	57	62	180	46	21	94	24	86	44	<1.0	<1.0	<1.0	3.4	<1.0	<1.0	<0.50	1.7
MW-712	UG/L	5/17/2012	8400	650	130	180	740	150	86	22	44	240	77	<1.0	<1.0	<1.0	3.0	<1.0	<1.0	<0.50	1.1
MW-712	UG/L	9/6/2012	10000	1100	27	47	110	40	110	97	49	88	33	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-712	UG/L	11/19/2012	670	55	5.8	8.1	37	8.6	5.9	<10	11	17	4.9	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-712	UG/L	2/12/2013	3200	690	75	100	460	76	130	<10	37	190	54	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-712	UG/L	4/12/2013	5800	540	56	93	390	68	180	<10	36	130	46	<1.0	<1.0	<1.0	2.4	<1.0	<1.0	<0.50	<1.0
MW-713	UG/L	2/9/2011	280	29	<0.50	<0.50	1.7	<0.50	3.5	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.2	<1.0	<1.0	<0.50	<1.0
MW-713	UG/L	4/8/2011	1000	150	<0.50	0.91	1.6	<0.50	75	120	2.8	<1.0	<1.0	<1.0	<1.0	<1.0	5.4	<1.0	<1.0	<0.50	<1.0
MW-713	UG/L	9/2/2011	310	73	3.0	1.7	7.8	3.6	71	100	11	7.0	1.6	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-713	UG/L	11/22/2011	3300	900	1.6	3.4	12	2.6	230	220	2.2	2.0	<1.0	<1.0	<1.0	<1.0	2.5	<1.0	<1.0	<0.50	<1.0
MW-713	UG/L	11/22/2011	3500	800	1.9	3.8	14	2.9	230	230	2.7	2.4	<1.0	<1.0	<1.0	<1.0	2.8	<1.0	<1.0	<0.50	<1.0
MW-713	UG/L	2/13/2012	5500	1900	2.2	4.6	9.8	2.5	390	160	<1.0	1.6	<1.0	<1.0	<1.0	<1.0	3.1	<1.0	<1.0	<0.50	<1.0
MW-713	UG/L	5/17/2012	5100	2300	2.3	5.3	6.0	1.3	400	110	3.6	1.1	<1.0	<1.0	<1.0	<1.0	2.2	<1.0	<1.0	<0.50	<1.0
MW-713	UG/L	9/6/2012	9600	1600	3.5	6.4	6.8	1.5	410	75	14	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-713	UG/L	11/19/2012	750	350	0.79	1.5	2.1	<0.50	190	73	13	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-713	UG/L	2/12/2013	5500	5300	7.0	16	33	1.0	720	<10	17	1.0	9.6	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-713	UG/L	4/11/2013	8200	4900	8.2	13	37	1.9	760	310	6.8	1.5	13	<1.0	<1.0	<1.0	2.0	<1.0	<1.0	<0.50	<1.0
MW-713	UG/L	4/11/2013	8300	5000	8.4	13	38	2.0	800	320	6.7	1.5	14	<1.0	<1.0	<1.0	2.1	<1.0	<1.0	<0.50	1 2
MW-714	UG/L	2/14/2011	370	1.3	<0.50	<0.50	<1.0	<0.50	10	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-714	UG/L	4/7/2011	16000	16	4.0	2.1	11	1.9	16	<10	23	4.7	1.4	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-714	UG/L	9/2/2011	500	3.8	<0.50	<0.50	1.1	<0.50	9.7	37	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-714	UG/L	11/22/2011	430	9.0	<0.50	<0.50	<1.0	<0.50	8.4	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-714	UG/L	11/22/2011	490	4.7	<0.50	<0.50	<1.0	<0.50	7.9	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-714	UG/L	2/13/2012	760	3.9	<0.50	<0.50	<1.0	<0.50	7.1	23	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-714	UG/L	2/13/2012	730	5.0	0.72	<0.50	1.1	<0.50	8.4	29	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-714	UG/L	5/18/2012	390	2.4	<0.50	<0.50	<1.0	<0.50	7.1	<10	1.2	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-714	UG/L	9/6/2012	500	1.6	<0.50	<0.50	<1.0	<0.50	2.3	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-714	UG/L	11/19/2012	<50	1.2	<0.50	<0.50	<1.0	<0.50	2.4	20	3.7	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-714	UG/L	2/12/2013	86	1.3	<0.50	<0.50	<1.0	<0.50	7.6	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-714	UG/L	4/11/2013	170	1.3	<0.50	<0.50	<1.0	<0.50	7.2	52	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-715	UG/L	2/14/2011	2000	480	12	1.7	24	7.4	2.8	<10	<1.0	2.6	4.2	<1.0	<1.0	<1.0	1.5	<1.0	<1.0	<0.50	<1.0
MW-715	UG/L	4/8/2011	1500	310	5.6	1.0	3.6	1.6	8.8	<10	3.8	<1.0	1.7	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-715	UG/L	9/2/2011	5500	800	2.5	4.0	12	5.3	8.2	22	5.0	4.5	4.8	<1.0	<1.0	<1.0	1.5	<1.0	<1.0	0.56	1.9
MW-715	UG/L	9/2/2011	1100	420	1.4	2.2	6.1	2.5	7.9	20	3.8	2.5	4.6	<1.0	<1.0	<1.0	1.5	<1.0	<1.0	0.53	1.2

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
MW-715	UG/L	11/22/2011	1500	450	1.5	6.0	<1.0	<0.50	8.5	11	3.5	4.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-715	UG/L	2/1/2012	860	270	2.6	1.7	5.6	1.1	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-715	UG/L	5/18/2012	13000	2100	19	1100	1900	350	4.3	<10	230	930	270	<1.0	<1.0	<1.0	1.4	<1.0	<1.0	<0.50	2.1
MW-715	UG/L	9/6/2012	610	11	0.56	62	<1.0	<0.50	1.2	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-715	UG/L	11/19/2012	<50	0.52	<0.50	<0.50	<1.0	<0.50	<1.0	<10	2.2	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-715	UG/L	2/12/2013	<50	0.71	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
MW-715	UG/L	4/12/2013	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-1	ug/L	11/1/1989		390	3.9	2.1								<0.5A		<0.5A			3.5A	<0.5A	21
W-1	ug/L	3/1/1990		140	<5	<5								<5		<10			<5	<5	<20
W-1	ug/L	4/1/1990		200	12	12								<5		<5	<25		1.6	<5	<5
W-1	ug/L	12/18/1996	800	78	<5	<5			<10		10	<5	<5	<5		<5	<5		<5	<5	<10
W-1	ug/L	1/14/1998	1100	62	<5	<5			<5		<10	<5	<5	<5		<5	<5		<5	<5	16
W-1	ug/L	8/20/1998	1200	79	<5	<5			14		<10	<5	<5	<5		<5	8.6		8.4	<5	26
W-1	ug/L	1/29/1999	1400	57	<5	<5			<5		<10	<5	<5	<5		<5	<5		<5	<5	18
W-1	ug/L	7/19/1999	1500	48	<2	<2			<2		<20	<2	<2	<2		<2	<2		<2	<1	<1
W-1	ug/L	8/3/2000	880	29	<1	<1			10		<10	<1	<1	<1		<1	1.6		1.6	<0.5	7.3
W-1	ug/L	2/8/2001	<500	21	<1	<1			68		<10	<1	<1	<1		<1	2.3		<1	<0.5	6.3
W-1	ug/L	7/26/2001	620	18	<1	<1			62		<10	<1	<1	<1		<1	2.8		1.8	<0.5	6.8
W-1	ug/L	5/8/2002	280	7.7	<1	<1			5.9	44000	<10	<1	<1	<1		<1	3.1		<1	<0.5	6.4
W-1	ug/L	9/25/2002	210	12	<1	<1			1.9	30000	<10	<1	<1	<1		<1	6.5		<1	<0.5	14
W-1	ug/L	7/1/2004	460	14	2.8	1.5	<0.5	<0.5	3J	<100	<5	<5	<5	<5		4J	9.3		1J	<5	2
W-1	ug/L	10/6/2005	310	43	<1	<1	<1	<1	25	34	<10	<1	<1	<1		1.6	<1		<1	<0.5	7.1
W-1	ug/L	2/15/2006	266	32	<5	<5	<5	<5	22	37	<5	<5	<5	<5		1.3	<5		<5	<5	3.3
W-1	ug/L	8/3/2006	1100	86	<2	<2	<2	<2	77	100	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-1	ug/L	11/9/2006	470	100	<2	<2	<2	<2	65	78	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-1	ug/L	2/8/2007	500	77	<2	<2	<2	<2	21	<50	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-1	ug/L	5/10/2007	890	110	0.57	0.61	<2	0.32	28	43	1	<2	<2	<2		0.42	<2		<2	<2	1.8
W-1	ug/L	8/9/2007	1100	140	0.84	0.84	<2	0.63	64	84	1.1	<2	<2	<2		0.47	<2		0.32	<2	1.9
W-1	ug/L	11/7/2007	1200	140	1.6	1.2	0.68	0.91	56	80	1.6	0.38	2.1	<0.32		0.7	<0.32		<0.27	<0.28	1.2
W-1	ug/L	2/7/2008	1000	96	<2	<2	<2	<2	31	51	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-1	ug/L	1/20/2009	230	15	<2	<2	<2	<2	3.1	23	<5	<2	<2	<2		0.87	<2		0.58	<2	2.8
W-1	ug/L	1/20/2009	220	19	<2	<2	<2	<2	3.9	35	<5	<2	<2	<2		1.1	0.4		0.61	<2	3.7
W-1	ug/L	4/24/2009	180	3.9	<2	<2	<2	<2	<5	26	<5	<2	<2	<2		1.4	<2		0.74	<2	9.5
W-1	UG/L	3/5/2010	270	3.3	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	1.3
W-1	UG/L	5/13/2010	260	9.3	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	1.2
W-1	UG/L	8/6/2010	260	17	<0.50	<0.50		<0.50	<1.0	10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-1	UG/L	11/5/2010	150	15	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-1	UG/L	2/4/2011	200	2.7	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-1	UG/L	4/14/2011	150	1.4	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-1	UG/L	8/26/2011	130	3.9	<0.50	<0.50	<1.0	<0.50	1.3	16	<1.0	<1.0	<1.0	<1.0	<1.0	4.2	<1.0	<1.0	<1.0	<0.50	6.4
W-1	UG/L	11/14/2011	160	12	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-1	UG/L	11/14/2011	160	12	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	5.1	<1.0	<1.0	<1.0	<0.50	<1.0
W-1	UG/L	2/6/2012	160	18	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	3.7	<1.0	<1.0	<1.0	<0.50	2.4
W-1	UG/L	5/7/2012	680	15	<0.50	<0.50	<1.0	<0.50	<1.0	23	<1.0	<1.0	<1.0	<1.0	<1.0	2.2	<1.0	<1.0	<1.0	<0.50	1.8
W-1	UG/L	8/27/2012	180	9.1	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-1	UG/L	11/5/2012	67	1.2	<0.50	<0.50	<1.0	<0.50	<1.0	<10	4.4	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-1	UG/L	1/30/2013	120	11	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
W-1	UG/L	4/3/2013	<50	1.2	<0.50	<0.50	<1.0	<0.50	<1.0	<10	5.6	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-10	ug/L	11/8/2006	26000	8200	5000	570	2100	820	<100	<1000	340	360	110	<40		<40	<40		<40	<40	<100
W-10	ug/L	2/9/2007	28000	6400	2200	520	2200	710	<500	<5000	<500	280	<200	<200		<200	<200		<200	<200	<500
W-10	ug/L	2/9/2007	26000	5100	1600	410	1800	570	<500	<5000	<500	260	<200	<200		<200	<200		<200	<200	<500
W-10	ug/L	5/11/2007	7900	430	140	100	480	130	<10	84	100	130	48	<4		<4	6		8.2	1.2	3 6
W-10	ug/L	5/11/2007	7800	500	160	110	540	150	<25	85	150	150	53	<10		<10	6.6		8.8	1.4	3 9
W-10	ug/L	8/9/2007	5400	590	20	82	330	40	<25	68	59	90	33	<10		<10	6.4		8	<10	3
W-10	ug/L	11/9/2007	<12000	4700	460	330	1300	240	<32	<490	240	190	55	<32		<27	<32		<27	<28	<30
W-10	ug/L	2/8/2008	<28000	7200	280	300	1300	190	<500	<5000	140	140	38	<200		<200	<200		<200	<200	<500
W-10	ug/L	2/8/2008	<25000	7600	310	330	1400	200	<500	<5000	170	150	42	<200		<200	<200		<200	<200	<500
W-10	ug/L	1/21/2009	20000	8100	<200	440	1400	<200	<500	<5000	<500	230	<200	<200		<200	<200		<200	<200	<500
W-10	ug/L	4/27/2009	16000	7400	<200	490	1400	<200	<500	<5000	270	230	36	<200		<200	<200		<200	<200	<500
W-10	ug/L	4/27/2009	15000	5100	<200	350	830	<200	<500	<5000	220	190	31	<200		<200	<200		<200	<200	<500
W-10	UG/L	3/8/2010	12000	4200	4.4	200		1.6	<1.0	<10	110	93	18	<1.0		<1.0	<1.0		<1.0	7.3	<1.0
W-10	UG/L	3/8/2010	8600	3100	<250	<250		<250	<500	<5000	<500	<500	<500	<500		<500	<500		<500	<250	<500
W-10	UG/L	5/17/2010	9500	3900	7.4	230		1.9	<1.0	<10	130	70	13	<1.0		<1.0	<1.0		<1.0	2.7	<1.0
W-10	UG/L	5/17/2010	10000	2900	10	160		1.7	<1.0	15	110	82	14	<1.0		<1.0	<1.0		<1.0	4.2	<1.0
W-10	UG/L	8/9/2010	7900	2400	12	130		1.9	<1.0	93	60	62	10	<1.0		<1.0	<1.0		<1.0	3.0	<1.0
W-10	UG/L	11/8/2010	7700	2900	45	160	140	6.4	<1.0	<10	180	56	8.1	<1.0	<1.0	<1.0	<1.0	<1.0	1.0	2.6	1.4
W-10	UG/L	2/8/2011	11000	2600	100	160	140	28	<1.0	<10	150	61	13	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	4.0	<1.0
W-10	UG/L	4/21/2011	12000	4900	97	240	190	38	<1.0	250	150	65	15	<1.0	<1.0	<1.0	<1.0	<1.0	1.6	12	<1.0
W-10	UG/L	9/1/2011	8200	2900	2.2	120	44	1.1	<1.0	140	97	31	5.7	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	4.9	<1.0
W-10	UG/L	11/16/2011	8800	840	3.9	190	92	1.1	<1.0	<10	94	49	10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-10	UG/L	2/8/2012	10000	3100	5.5	230	150	2.9	<1.0	<10	130	73	12	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	5.6	<1.0
W-10	UG/L	5/10/2012	1000	15	<0.50	1.4	1.2	<0.50	<1.0	<10	21	4.3	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-10	UG/L	8/28/2012	8200	3100	4.3	160	32	1.4	<1.0	61	270	27	2.8	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.8	<1.0
W-10	UG/L	11/7/2012	5100	930	7.9	120	65	2.9	<1.0	65	130	27	4.2	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.3	<1.0
W-10	UG/L	1/29/2013	160	4.4	8.1	5.6	22	9.9	<1.0	35	71	15	3.3	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-10	UG/L	4/1/2013	490	6.3	<0.50	<0.50	<1.0	<0.50	<1.0	150	13	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-11	ug/L	11/9/2006	5200	99	12	74	240	37	<5	<50	<5	73	40	<2		<2	18		<2	<2	<5
W-11	ug/L	11/9/2006	12000	96	7.8	54	140	21	<5	<50	<5	60	34	<2		<2	18		<2	<2	<5
W-11	ug/L	2/9/2007	8000	95	14	78	280	27	<10	<100	<10	56	28	<4		<4	15		<4	<4	<10
W-11	ug/L	5/9/2007	540	45	1.6	19	47	3.1	<5	<50	0.68	9	4.4	<2		0.41	18		<2	<2	0.96
W-11	ug/L	8/8/2007	<1100	700	3.7	36	11	7.1	<5	<50	0.81	15	8.6	<2		<2	9.9		<2	0.29	1.1
W-11	ug/L	11/8/2007	460	61	1.2	14	37	13	<0.32	<4.9	1	35	17	<0.32		<0.27	10		<0.27	<0.28	<0.3
W-11	UG/L	12/8/2010	77000	150	51	260	2300	690	17	43	48	1300	800	<1.0	<1.0	<1.0	<1.0	<1.0	1.4	<0.50	<1.0
W-11	UG/L	2/4/2011	10000	100	1.2	23	100	16	<1.0	<10	7.6	100	180	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-11	UG/L	4/15/2011	6300	410	15	50	390	18	<1.0	<10	3.4	83	280	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-11	UG/L	8/29/2011	10000	560	2.2	57	640	14	<1.0	<10	<1.0	100	190	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-11	UG/L	11/14/2011	10000	620	3.0	100	510	7.5	<1.0	<10	6.0	130	240	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-11	UG/L	2/8/2012	2900	12	<0.50	6.2	50	0.80	<1.0	<10	2.7	24	39	<1.0	<1.0	<1.0	2.0	<1.0	<1.0	0.90	<1.0
W-11	UG/L	5/10/2012	1800	8.4	<0.50	3.1	7.3	0.80	<1.0	<10	1.7	4.6	10	<1.0	<1.0	<1.0	2.0	<1.0	<1.0	0.50	<1.0
W-11	UG/L	8/28/2012	7400	16	30	47	130	20	<1.0	<10	5.0	70	97	<1.0	<1.0	<1.0	2.1	<1.0	<1.0	<0.50	<1.0
W-11	UG/L	11/8/2012	340	23	3.1	1.6	23	2.0	<1.0	<10	2.5	5.0	63	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-11	UG/L	11/19/2012	1400	24	1.6	0.82	6.2	<0.50	<1.0	<10	3.0	3.1	60	<1.0	<1.0	<1.0	5.3	<1.0	<1.0	<0.50	1.3
W-11	UG/L	1/31/2013	300	25	8.2	3.7	25	8.9	<1.0	<10	33	14	42	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
W-11	UG/L	4/5/2013	250	14	0.75	1.2	3.2	0.57	<1.0	<10	<1.0	1.4	8.9	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-12	ug/L	11/8/2006	1400	<2	<2	<2	<2	<2	<5	55	<5	<2	<2	<2		<2	5.4		<2	<2	<5
W-12	ug/L	2/7/2007	4800	<2	<2	<2	<2	<2	<5	50	<5	<2	<2	<2		<2	6.8		<2	<2	<5
W-12	ug/L	5/9/2007	220	<2	<2	<2	<2	<2	<5	40	<5	<2	<2	<2		0.31	4.3		<2	0.37	1.1
W-12	ug/L	8/8/2007	1100	<2	<2	0.56	<2	<2	0.36	40	<5	<2	<2	<2		<2	3.1		<2	<2	0.85
W-12	ug/L	11/6/2007	1500	0.37	<0.36	0.97	<0.6	<0.3	1.2	58	0.66	<0.23	<0.26	<0.32		<0.27	2.6		<0.27	0.42	0.47
W-12	ug/L	2/8/2008	410	0.94	<2	3	<2	<2	0.82	54	2.5	<2	<2	<2		<2	1.8		<2	0.45	<5
W-12	ug/L	1/20/2009	620	<2	<2	0.69	<2	<2	<5	32	<5	<2	<2	<2		0.48	5.4		<2	<2	2.4
W-12	ug/L	4/22/2009	1100	<2	<2	2.1	<2	<2	0.33	30	8.2	0.26	<2	<2		<2	3.7		<2	<2	1.5
W-12	UG/L	3/4/2010	400	<0.50	<0.50	2.1		<0.50	<1.0	<10	1.5	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-12	UG/L	5/12/2010	610	<0.50	<0.50	3.0		<0.50	<1.0	<10	2.1	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-12	UG/L	8/5/2010	650	<0.50	<0.50	3.5		<0.50	<1.0	<10	2.8	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-12	UG/L	11/4/2010	530	<0.50	<0.50	1.4	<1.0	<0.50	<1.0	<10	1.7	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-12	UG/L	2/3/2011	310	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-12	UG/L	4/19/2011	220	<0.50	<0.50	0.57	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.1	<1.0	<1.0	<0.50	2.7
W-12	UG/L	8/25/2011	360	<0.50	<0.50	1.3	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-12	UG/L	11/14/2011	63	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.2	<1.0	<1.0	<0.50	<1.0
W-12	UG/L	2/8/2012	400	<0.50	<0.50	2.2	<1.0	<0.50	<1.0	<10	1.6	<1.0	<1.0	<1.0	<1.0	<1.0	2.3	<1.0	<1.0	<0.50	2.2
W-12	UG/L	5/9/2012	450	<0.50	<0.50	0.59	<1.0	<0.50	<1.0	27	1.2	<1.0	<1.0	<1.0	<1.0	<1.0	1.4	<1.0	<1.0	<0.50	1.2
W-12	UG/L	8/30/2012	580	<0.50	<0.50	1.5	1.0	<0.50	<1.0	<10	20	1.2	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-12	UG/L	11/8/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-12	UG/L	1/31/2013	<50	<0.50	1.2	1.6	8.4	4.0	<1.0	<10	34	10	2.9	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-12	UG/L	4/2/2013	150	<0.50	<0.50	3.0	17	4.2	<1.0	26	13	15	4.6	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-14A	ug/L	2/12/2008	42	<2	<2	<2	<2	<2	<5	<50	<5	<2	<2	2.3		1.1	9		0.46	0.37	<5
W-14A	ug/L	1/13/2009	<50	<2	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-14A	ug/L	4/21/2009	54	<2	<2	<2	<2	<2	0.47	8.1	<5	<2	<2	1.3		0.86	8.7		0.44	0.4	<5
W-14A	UG/L	3/1/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	1.7		<1.0	<0.50	<1.0
W-14A	UG/L	5/10/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	1.9		<1.0	<0.50	<1.0
W-14A	UG/L	8/2/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	3.4		<1.0	<0.50	<1.0
W-14A	UG/L	11/1/2010	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-14A	UG/L	1/31/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-14A	UG/L	4/4/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.6	<1.0	<1.0	<0.50	<1.0
W-14A	UG/L	8/22/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	5.8	1.0	5.2	<1.0	<1.0	<0.50	<1.0
W-14A	UG/L	11/7/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.8	<1.0	<1.0	<0.50	<1.0
W-14A	UG/L	1/30/2012	200	1.5	<0.50	38	<1.0	<0.50	<1.0	<10	<1.0	1.1	<1.0	<1.0	3.2	<1.0	10	1.4	<1.0	<0.50	<1.0
W-14A	UG/L	5/1/2012	390	41	<0.50	9.5	1.3	2.7	2.9	<10	<1.0	1.2	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-14A	UG/L	8/20/2012	1600	500	16	34	78	64	2.9	<10	110	57	20	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-14A	UG/L	10/26/2012	3800	4500	5.1	150	240	110	1.5	<10	51	120	42	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	4.4
W-14A	UG/L	1/22/2013	1100	110	<0.50	33	2.2	<0.50	<1.0	<10	13	13	<1.0	<1.0	2.5	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-14A	UG/L	4/1/2013	96	5.8	1.8	1.4	6.6	4.5	<1.0	<10	91	12	2.4	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-14B	ug/L	2/12/2008	<50	<2	<2	<2	<2	<2	<5	<50	<5	<2	<2	0.72		<2	0.83		<2	<2	<5
W-14B	ug/L	1/13/2009	170	<2	<2	<2	<2	<2	<5	<50	<5	<2	<2	8.4		<2	4.8		<2	<2	<5
W-14B	ug/L	4/21/2009	65	<2	<2	<2	<2	<2	<5	<50	<5	<2	<2	19		2.6	9.6		2.2	0.45	<5
W-14B	UG/L	3/1/2010	99	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	5.6		<1.0	<0.50	<1.0
W-14B	UG/L	5/10/2010	99	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	1.2		1.1	6.2		<1.0	<0.50	<1.0

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
W-14B	UG/L	8/2/2010	55	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	3.1		<1.0	<0.50	<1.0
W-14B	UG/L	11/1/2010	88	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	2.0	45	2.0	10	14	1.2	<0.50	<1.0
W-14B	UG/L	1/31/2011	65	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	9.7	<1.0	2.0	3.1	<1.0	<0.50	<1.0
W-14B	UG/L	4/4/2011	<50	<0.50	1.8	<0.50	<1.0	<0.50	<1.0	48	<1.0	<1.0	<1.0	15	99	2.8	13	34	2.9	0.53	<1.0
W-14B	UG/L	8/22/2011	200	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	28	130	2.4	9.8	53	3.2	0.98	<1.0
W-14B	UG/L	11/7/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	5.1	<1.0	<1.0	1.8	<1.0	<0.50	<1.0
W-14B	UG/L	1/30/2012	220	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	22	100	<1.0	12	55	3.1	<0.50	<1.0
W-14B	UG/L	5/1/2012	150	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	69	<1.0	<1.0	<1.0	8.0	82	<1.0	11	53	2.4	<0.50	<1.0
W-14B	UG/L	8/20/2012	180	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	56	<1.0	<1.0	<1.0	8.9	150	2.4	13	60	2.9	<0.50	<1.0
W-14B	UG/L	10/26/2012	52	6.0	<0.50	1.6	4.8	0.89	<1.0	<10	20	1.8	<1.0	4.3	82	1.6	7.4	31	1.6	<0.50	1.9
W-14B	UG/L	1/22/2013	150	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	7.4	98	<1.0	14	44	2.0	<0.50	<1.0
W-14B	UG/L	4/1/2013	<50	<0.50	0.61	<0.50	2.0	1.2	<1.0	<10	21	2.6	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-14C	ug/L	2/12/2008	260	1.2	<2	<2	<2	<2	<5	<50	<5	<2	<2	0.89		5.7	22		3.7	0.48	0.58
W-14C	ug/L	1/14/2009	120	2.5	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		8.8	34		3.4	<2	<5
W-14C	ug/L	4/21/2009	67	1.5	<2	<2	<2	<2	<5	10	<5	<2	<2	<2		4.5	23		2.1	<2	<5
W-14C	UG/L	3/1/2010	300	1.6	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		5.8	34		2.4	<0.50	<1.0
W-14C	UG/L	5/10/2010	120	0.58	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		2.0	13		<1.0	<0.50	<1.0
W-14C	UG/L	8/2/2010	77	1.1	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		4.6	35		2.4	<0.50	<1.0
W-14C	UG/L	11/1/2010	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-14C	UG/L	1/31/2011	60	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	3.8	1.1	9.9	3.0	<1.0	<0.50	<1.0
W-14C	UG/L	4/4/2011	<50	1.2	<0.50	<0.50	<1.0	<0.50	<1.0	27	<1.0	<1.0	<1.0	<1.0	24	3.9	30	16	3.1	<0.50	<1.0
W-14C	UG/L	8/22/2011	290	0.73	<0.50	<0.50	<1.0	<0.50	<1.0	22	<1.0	<1.0	<1.0	<1.0	21	2.3	26	12	2.2	<0.50	<1.0
W-14C	UG/L	11/7/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	1.2	<1.0	3.2	<1.0	<1.0	<0.50	<1.0
W-14C	UG/L	1/30/2012	100	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	3.4	<1.0	5.3	2.2	<1.0	<0.50	<1.0
W-14C	UG/L	5/1/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	3.8	<1.0	<1.0	<0.50	<1.0
W-14C	UG/L	8/20/2012	71	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	5.8	1.4	<1.0	<0.50	<1.0
W-14C	UG/L	10/26/2012	<50	0.75	<0.50	<0.50	<1.0	<0.50	<1.0	<10	6.1	<1.0	<1.0	<1.0	<1.0	<1.0	8.4	2.6	<1.0	<0.50	2.6
W-14C	UG/L	1/22/2013	110	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	7.5	2.6	1.1	<0.50	<1.0
W-14C	UG/L	4/1/2013	<50	<0.50	0.88	0.58	2.7	1.7	<1.0	<10	27	2.4	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15A	ug/L	2/11/2008	2700	620	4.9	5.1	11	<20	650	120	<50	<20	<20	<20		<20	<20		<20	<20	<50
W-15A	ug/L	1/14/2009	230	7.4	<2	<2	<2	<2	190	170	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-15A	ug/L	4/24/2009	530	8.4	<4	<4	<4	<4	220	220	<10	<4	<4	<4		<4	<4		<4	<4	<10
W-15A	UG/L	3/2/2010	240	0.93	<0.50	<0.50		<0.50	44	94	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-15A	UG/L	5/10/2010	260	1.5	<0.50	<0.50		<0.50	85	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-15A	UG/L	8/2/2010	310	0.54	<0.50	<0.50		<0.50	71	180	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-15A	UG/L	11/1/2010	61	<0.50	<0.50	<0.50	<1.0	<0.50	2.5	88	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15A	UG/L	11/1/2010	74	0.66	<0.50	<0.50	1.0	<0.50	6.8	98	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15A	UG/L	2/1/2011	14000	1400	610	400	1800	400	260	390	64	490	200	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.6	<1.0
W-15A	UG/L	4/5/2011	22000	<0.50	<0.50	<0.50	<1.0	<0.50	450	<10	150	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15A	UG/L	2/2/2012	62000	4400	2400	2400	9900	2300	930	<10	4.6	2900	880	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15A	UG/L	5/2/2012	2100000	3900	3600	3900	13000	4400	940	220	450	6200	1800	<10	<10	<10	<10	<10	<10	<5.0	<10
W-15A	UG/L	8/21/2012	23000	540	370	590	3300	620	160	<250	190	1100	340	<25	<25	<25	<25	<25	<25	<12	<25
W-15A	UG/L	10/30/2012	4500	41	23	46	260	75	39	120	330	270	120	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15A	UG/L	1/23/2013	2400	100	36	57	200	95	57	<10	120	170	94	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15A	UG/L	4/2/2013	3400	86	32	79	460	130	72	120	260	230	67	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
W-15B	ug/L	2/11/2008	<1600	900	<20	<20	7	<20	20	110	<50	<20	<20	<20		<20	<20		<20	<20	<50
W-15B	ug/L	1/14/2009	340	160	<2	<2	5	<2	20	110	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-15B	ug/L	4/24/2009	63	6.2	<2	<2	<2	<2	5.8	98	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-15B	UG/L	3/2/2010	220	3.8	<0.50	<0.50		<0.50	5.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-15B	UG/L	5/11/2010	230	20	<0.50	<0.50		<0.50	17	36	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-15B	UG/L	8/3/2010	250	14	<0.50	<0.50		<0.50	19	67	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-15B	UG/L	11/2/2010	740	38	<0.50	<0.50	3.2	0.74	50	87	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15B	UG/L	2/1/2011	120	7.0	1.7	0.55	4.0	1.4	22	21	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15B	UG/L	4/5/2011	1500	<0.50	66	18	120	64	130	<10	6.3	16	16	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15B	UG/L	8/23/2011	1400	120	40	17	110	30	260	210	<1.0	13	7.2	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15B	UG/L	8/23/2011	1100	110	34	15	100	29	200	220	<1.0	14	7.2	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15B	UG/L	11/10/2011	250	17	5.4	2.8	17	3.9	55	<10	<1.0	2.4	1.1	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15B	UG/L	2/2/2012	280	35	14	4.4	31	18	100	80	<1.0	2.3	3.8	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15B	UG/L	5/2/2012	780	27	2.6	3.1	18	6.3	200	160	<1.0	4.4	2.6	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15B	UG/L	8/20/2012	98	2.6	<0.50	<0.50	<1.0	0.52	110	87	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15B	UG/L	10/30/2012	190	9.2	2.2	1.5	12	2.7	49	96	43	4.0	1.4	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15B	UG/L	1/23/2013	300	3.8	1.9	9.0	65	15	12	<10	38	71	25	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15B	UG/L	4/2/2013	430	1.0	2.3	13	87	19	19	180	46	62	19	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15C	ug/L	2/11/2008	<50	0.94	0.57	<2	<2	<2	<5	18	<5	<2	<2	<2		<2	1.1		0.45	0.35	0.34
W-15C	ug/L	1/15/2009	29	1.1	<2	<2	<2	<2	<5	27	<5	<2	<2	<2		<2	5.7		1.2	0.86	0 9
W-15C	ug/L	4/24/2009	43	<2	<2	<2	<2	<2	<5	25	<5	<2	<2	<2		<2	1		<2	<2	<5
W-15C	UG/L	3/2/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	1.4		<1.0	<0.50	<1.0
W-15C	UG/L	5/11/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	1.6		<1.0	<0.50	<1.0
W-15C	UG/L	8/3/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	20	<1.0	<1.0	<1.0	<1.0		<1.0	4.7		1.0	0.54	1.5
W-15C	UG/L	11/2/2010	70	<0.50	<0.50	<0.50	<1.0	<0.50	2.9	<10	<1.0	<1.0	<1.0	<1.0	1.0	<1.0	1.7	<1.0	<1.0	<0.50	<1.0
W-15C	UG/L	2/1/2011	94	1.6	0.85	<0.50	2.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	1.8	<1.0	2.6	<1.0	<1.0	<0.50	<1.0
W-15C	UG/L	4/5/2011	120	10	4.8	1.9	10	2.6	4.2	<10	1.1	<1.0	<1.0	<1.0	4.6	<1.0	6.6	1.5	1.4	<0.50	1.8
W-15C	UG/L	8/23/2011	89	9.5	3.5	1.4	13	2.7	5.2	<10	<1.0	1.8	<1.0	<1.0	5.5	<1.0	6.5	1.6	<1.0	<0.50	<1.0
W-15C	UG/L	11/8/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15C	UG/L	1/31/2012	53	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	10	<1.0	<1.0	<1.0	<1.0	4.9	<1.0	5.8	1.5	<1.0	<0.50	<1.0
W-15C	UG/L	5/2/2012	60	0.64	0.67	1.4	6.4	1.3	<1.0	<10	<1.0	3.2	1.2	<1.0	1.3	<1.0	2.1	<1.0	<1.0	<0.50	<1.0
W-15C	UG/L	8/21/2012	140	4.1	1.7	0.92	5.9	1.4	1.7	10	2.9	1.5	<1.0	<1.0	3.7	<1.0	5.2	1.2	<1.0	<0.50	<1.0
W-15C	UG/L	10/30/2012	120	16	4.9	3.2	36	7.1	3.4	<10	9.9	6.6	2.9	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15C	UG/L	1/23/2013	180	<0.50	0.80	4.4	33	7.2	<1.0	<10	19	43	15	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-15C	UG/L	4/2/2013	410	20	5.8	9.8	86	21	6.3	25	30	42	13	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-16A	ug/L	11/9/2007	260	41	<0.36	<0.25	<0.6	<0.3	<0.32	30	<0.41	<0.23	<0.26	<0.32		<0.27	<0.32		2.6	<0.28	16
W-16A	ug/L	2/6/2008	310	40	<2	<2	<2	<2	<5	34	<5	<2	0.63	<2		0.88	<2		2.8	<2	14
W-16A	ug/L	1/21/2009	290	30	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		<2	<2		2.5	<2	7 2
W-16A	ug/L	4/27/2009	410	34	<2	<2	<2	<2	<5	20	<5	<2	0.27	<2		0.54	<2		1.8	<2	17
W-16A	UG/L	3/5/2010	220	4.2	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	2.9
W-16A	UG/L	5/14/2010	110	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-16A	UG/L	8/9/2010	120	0.93	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-16A	UG/L	11/5/2010	90	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-16A	UG/L	2/7/2011	320	12	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.7	<0.50	1.1
W-16A	UG/L	4/18/2011	520	24	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.2	<0.50	2.2
W-16A	UG/L	8/26/2011	280	13	<0.50	<0.50	<1.0	<0.50	<1.0	30	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.2	<0.50	<1.0

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
W-16A	UG/L	11/8/2011	65	3.1	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-16A	UG/L	2/3/2012	230	16	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.1	<0.50	<1.0
W-16A	UG/L	5/3/2012	550	22	<0.50	1.0	4.4	1.1	<1.0	<10	<1.0	1.8	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.2	<0.50	<1.0
W-16A	UG/L	8/22/2012	390	11	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-16A	UG/L	10/31/2012	86	6.9	<0.50	<0.50	<1.0	<0.50	<1.0	<10	3.9	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-16A	UG/L	1/23/2013	84	<0.50	<0.50	0.81	5.4	1.3	<1.0	43	6.5	11	3.4	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-16A	UG/L	4/3/2013	340	20	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.0	<0.50	<1.0
W-16B	ug/L	11/9/2007	37	7.4	<0.36	<0.25	<0.6	<0.3	<0.32	9.1	0.8	0.26	<0.26	<0.32		8.7	6.6		<0.27	<0.28	<0.3
W-16B	ug/L	2/6/2008	400	48	<2	0.33	<2	<2	<5	9.9	1.9	0.4	<2	<2		43	27		<2	<2	<5
W-16B	ug/L	1/21/2009	73	16	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		15	9.7		<2	<2	<5
W-16B	ug/L	4/27/2009	47	0.9	<20	<20	<20	<20	<50	<500	<50	<20	<20	<20		9.4	6.1		<20	<20	<50
W-16B	UG/L	3/8/2010	73	8.6	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		3.7	5.8		<1.0	<0.50	<1.0
W-16B	UG/L	5/14/2010	60	3.0	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		1.0	3.0		<1.0	<0.50	<1.0
W-16B	UG/L	8/9/2010	<50	1.3	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-16B	UG/L	11/5/2010	110	23	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	9.4	13	<1.0	1.2	<0.50	<1.0
W-16B	UG/L	2/7/2011	290	80	<0.50	<0.50	<1.0	<0.50	<1.0	<10	18	<1.0	<1.0	<1.0	3.5	50	70	2.0	8.5	<0.50	2.9
W-16B	UG/L	4/18/2011	550	100	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	6.4	100	89	2.6	9.2	<0.50	10
W-16B	UG/L	8/26/2011	89	20	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	12	16	<1.0	1.4	<0.50	1.1
W-16B	UG/L	11/8/2011	<50	24	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	1.0	19	13	<1.0	1.5	<0.50	<1.0
W-16B	UG/L	2/3/2012	210	30	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	1.4	24	16	<1.0	1.3	<0.50	<1.0
W-16B	UG/L	5/3/2012	410	150	<0.50	0.58	2.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	3.2	100	52	1.2	6.8	<0.50	23
W-16B	UG/L	8/22/2012	61	8.7	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	3.5	6.0	<1.0	<1.0	<0.50	<1.0
W-16B	UG/L	10/31/2012	58	13	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	6.6	4.2	<1.0	<1.0	<0.50	15
W-16B	UG/L	1/23/2013	80	3.5	<0.50	<0.50	2.9	0.82	<1.0	<10	4.9	5.7	1.8	<1.0	<1.0	<1.0	1.7	<1.0	<1.0	<0.50	<1.0
W-16B	UG/L	4/3/2013	<50	3.4	<0.50	<0.50	1.1	0.79	<1.0	<10	9.1	1.5	<1.0	<1.0	<1.0	1.2	1.9	<1.0	<1.0	<0.50	<1.0
W-16C	ug/L	11/9/2007	170	18	<0.36	<0.25	<0.6	<0.3	<0.32	13	<0.41	<0.23	<0.26	<0.32		12	40		11	<0.28	5 6
W-16C	ug/L	2/6/2008	360	30	0.46	<2	<2	<2	<5	21	<5	<2	<2	<2		14	66		24	<2	18
W-16C	ug/L	1/21/2009	510	40	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		17	73		35	<2	24
W-16C	ug/L	4/28/2009	170	20	<2	<2	<2	<2	<5	8.2	<5	<2	<2	<2		12	41		14	<2	8 2
W-16C	UG/L	3/8/2010	95	2.5	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		1.2	9.1		1.6	<0.50	<1.0
W-16C	UG/L	5/14/2010	63	1.3	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	3.8		1.2	<0.50	<1.0
W-16C	UG/L	8/9/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-16C	UG/L	8/9/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-16C	UG/L	11/5/2010	390	14	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	7.6	30	1.4	11	<0.50	9.6
W-16C	UG/L	2/7/2011	440	33	0.54	<0.50	<1.0	<0.50	<1.0	<10	6.9	<1.0	<1.0	<1.0	<1.0	15	68	3.3	22	<0.50	14
W-16C	UG/L	4/18/2011	510	39	0.51	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	1.2	20	80	4.7	32	<0.50	30
W-16C	UG/L	8/26/2011	320	30	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	15	63	2.8	24	<0.50	16
W-16C	UG/L	11/9/2011	270	24	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	1.2	16	58	2.1	16	<0.50	<1.0
W-16C	UG/L	2/3/2012	250	23	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	1.0	16	54	2.8	17	<0.50	<1.0
W-16C	UG/L	5/3/2012	380	14	<0.50	<0.50	2.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	10	32	<1.0	9.8	<0.50	10
W-16C	UG/L	8/22/2012	520	22	<12	<12	<25	<12	<25	<250	<25	<25	<25	<25	<25	<25	42	<25	<25	<12	<25
W-16C	UG/L	10/31/2012	140	10	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	6.7	16	<1.0	8.0	<0.50	50
W-16C	UG/L	1/23/2013	58	<0.50	<0.50	<0.50	2.8	0.81	<1.0	<10	6.0	4.9	1.6	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-16C	UG/L	4/3/2013	<50	0.70	<0.50	<0.50	<1.0	0.56	<1.0	<10	7.8	<1.0	<1.0	<1.0	<1.0	<1.0	2.6	<1.0	<1.0	<0.50	<1.0
W-17A	ug/L	2/14/2008	100	<2	<2	<2	<2	<2	<5	140	<5	<2	<2	<2		<2	6.2		0.47	1.4	0.7

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
W-17A	ug/L	1/16/2009	78	<2	<2	<2	<2	<2	<5	54	0.41	0.33	<2	<2		0.39	1.4		<2	<2	<5
W-17A	ug/L	4/22/2009	180	4.5	<2	<2	<2	<2	<5	57	<5	<2	<2	<2		1.9	7.7		0.51	0.65	<5
W-17A	UG/L	3/3/2010	51	<0.50	<0.50	<0.50		<0.50	<1.0	14	<1.0	<1.0	<1.0	<1.0		<1.0	1.6		<1.0	<0.50	<1.0
W-17A	UG/L	5/12/2010	110	1.1	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	4.2		<1.0	<0.50	<1.0
W-17A	UG/L	8/4/2010	56	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	1.7		<1.0	<0.50	<1.0
W-17A	UG/L	11/3/2010	69	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.3	<1.0	<1.0	<0.50	<1.0
W-17A	UG/L	2/2/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.0	<1.0	<1.0	<0.50	<1.0
W-17A	UG/L	4/20/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	38	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.9	<1.0	<1.0	<0.50	<1.0
W-17A	UG/L	8/24/2011	98	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.5	<1.0	<1.0	<0.50	<1.0
W-17A	UG/L	11/9/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	9.6	<1.0	<1.0	<0.50	<1.0
W-17A	UG/L	2/7/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	17	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.5	<1.0	<1.0	<0.50	<1.0
W-17A	UG/L	5/4/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	3.8	<1.0	<1.0	<0.50	<1.0
W-17A	UG/L	8/23/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	12	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.9	<1.0	<1.0	<0.50	<1.0
W-17A	UG/L	11/1/2012	100	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	12	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	6.6	1.1	<1.0	<0.50	<1.0
W-17A	UG/L	1/25/2013	<50	<0.50	<0.50	<0.50	1.7	<0.50	<1.0	<10	2.6	2.4	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17A	UG/L	4/9/2013	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17B	ug/L	2/14/2008	39	<2	<2	<2	<2	<2	<5	30	<5	<2	<2	<2		<2	1.4		<2	<2	<5
W-17B	ug/L	1/16/2009	38	<2	<2	<2	<2	<2	<5	18	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-17B	ug/L	4/22/2009	<50	<2	<2	<2	<2	<2	<5	18	<5	<2	<2	<2		<2	0.71		<2	<2	<5
W-17B	UG/L	3/3/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-17B	UG/L	5/12/2010	54	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-17B	UG/L	8/5/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-17B	UG/L	11/3/2010	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17B	UG/L	2/2/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17B	UG/L	4/20/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	35	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17B	UG/L	8/24/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17B	UG/L	11/9/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17B	UG/L	2/7/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	14	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17B	UG/L	5/4/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17B	UG/L	8/23/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17B	UG/L	11/1/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	24	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17B	UG/L	1/25/2013	<50	<0.50	<0.50	<0.50	2.0	0.64	<1.0	<10	3.1	2.6	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17B	UG/L	4/9/2013	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17C	ug/L	2/14/2008	36	<2	<2	<2	<2	<2	<5	25	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-17C	ug/L	1/16/2009	29	<2	<2	<2	<2	<2	<5	21	<5	<2	<2	<2		<2	1.2		<2	<2	<5
W-17C	ug/L	4/23/2009	<50	<2	<2	<2	<2	<2	<5	18	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-17C	UG/L	3/4/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-17C	UG/L	5/12/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-17C	UG/L	8/5/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-17C	UG/L	11/3/2010	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17C	UG/L	2/2/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17C	UG/L	4/20/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	31	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17C	UG/L	8/24/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17C	UG/L	11/9/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17C	UG/L	2/7/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17C	UG/L	5/4/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	11	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
W-17C	UG/L	8/23/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17C	UG/L	11/1/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	11	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17C	UG/L	1/25/2013	<50	<0.50	<0.50	<0.50	1.3	<0.50	<1.0	<10	2.0	1.7	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-17C	UG/L	4/9/2013	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-3A	ug/L	1/13/1998	4300000	150000	<6000	35000			<200000												
W-3A	ug/L	8/20/1998	1100	220	<25	33			440		350	<25	<25	<25		<25	<25		<25	<25	<50
W-3A	ug/L	1/28/1999	690	160	<50	<50			340		240	<50	<50	<50		<50	<50		<50	<50	<100
W-3A	ug/L	7/19/1999	5400	120	<20	<20			380		<200	37	<20	<20		<20	<20		<20	<10	<10
W-3A	ug/L	1/13/2000	14000	140	<10	<10			210		<100	<10	<10	<10		<10	<10		<10	<5	7
W-3A	ug/L	8/4/2000	3400	170	<20	8.4			220		<50	2	2	<2		<2	<20		<20	<1	5
W-3A	ug/L	2/8/2001	2700	34	<1	2.9			12		63	13	4.4	<1		<1	<1		<1	<0.5	1.7
W-3A	ug/L	7/26/2001	3400	42	<1	1.7			6.2		11	15	<1	<1		<1	<1		<1	<0.5	27
W-3A	ug/L	5/6/2002	NS	NS	NS	NS			NS	NS	NS	NS	NS	NS		NS	NS		NS	NS	NS
W-3A	ug/L	9/25/2002	NS	NS	NS	NS			NS	NS	NS	NS	NS	NS		NS	NS		NS	NS	NS
W-3A	ug/L	2/16/2006	306	<1	<5	<5	<5	<5	6.2	16	<5	18	16	<5		<5	<5		<5	<5	<5
W-3A	ug/L	8/3/2006	39000	<2	<2	<2	<2	<2	9	<50	38	<2	<2	<2		<2	<2		<2	<2	<5
W-3A	ug/L	11/9/2006	8100	<2	<2	<2	<2	<2	11	<50	37	6.4	9.5	<2		<2	<2		<2	<2	<5
W-3A	ug/L	2/8/2007	1400	<2	<2	<2	<2	<2	8.4	<50	30	3.9	6.1	<2		<2	<2		<2	<2	<5
W-3A	ug/L	5/10/2007	14000	0.66	<2	<2	<2	<2	7.8	23	16	2.3	3.6	<2		<2	<2		<2	<2	<5
W-3A	ug/L	8/9/2007	1900	0.79	<2	<2	<2	0.34	9.8	26	14	2	2.3	<2		<2	<2		<2	<2	<5
W-3A	ug/L	11/7/2007	1500	0.62	<0.36	<0.25	<0.6	<0.3	9.7	26	<0.41	0.64	0.67	<0.32		<0.27	<0.32		<0.27	<0.28	<0.3
W-3A	ug/L	2/7/2008	180	<2	<2	<2	<2	<2	10	<50	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-4	ug/L	3/1/1990		120	<0.5	19								<0.5		<0.5	3.2		8.3	<0.5	<0.5
W-4	ug/L	4/1/1990		28	1.4	4.8								<1		<1	0.81		2.2	<1	4.3
W-4	ug/L	12/18/1996	420	80	<5	<5			<10		<5	<5	<5	<5		<5	<5		<5	<5	<10
W-4	ug/L	1/14/1998	920	120	<5	<5			<5		<10	<5	<5	<5		<5	<5		<5	<5	16
W-4	ug/L	8/20/1998	500	57	<5	<5			18		<10	<5	<5	<5		<5	<5		<5	<5	9 8
W-4	ug/L	1/29/1999	460	55	<5	<5			20		<10	<5	<5	<5		<5	<5		<5	<5	11
W-4	ug/L	7/19/1999	710	72	<2	<2			<2		<20	<2	<2	<2		<2	<2		<2	<1	<1
W-4	ug/L	1/13/2000	660	49	<1	<1			<1		<10	<1	<1	<1		<1	1.3		<1	<0.5	13
W-4	ug/L	8/3/2000	<500	47	<1	<1					<10	<1	<1	<1		1 2	<1		<1	<0.5	12
W-4	ug/L	2/8/2001	<500	42	<1	<1			<1		<10	<1	<1	<1		<1	<1		1.1	0.67	7
W-4	ug/L	7/26/2001	320	42	<1	<1			<1		<10	<1	<1	<1		<1	<1		1	<0.5	<0.5
W-4	ug/L	5/8/2002	250	33	<1	<1			<1	60000	<10	<1	<1	<1		2	<1		1.3	<0.5	5.2
W-4	ug/L	9/25/2002	290	62	<1	<1			<1	45000	<1	<1	<1	<1		3.8	<1		2	<0.5	<0.5
W-4	ug/L	7/1/2004	350	30	2.6	1.9	0.66	<0.5	<5	<100	<5	<5	<5	<5		1J	3J		2J	<5	11
W-4	ug/L	10/6/2005	350	31	<1	<1	<1	<1	<1	47	<10	<1	<1	<1		<1	6.4		1.7	<0.5	1.3
W-4	ug/L	2/15/2006	501	43	<5	<5	<5	<5	<1	38	<5	<5	<5	<5		<5	2.8		2.5	<5	2.4
W-4	ug/L	8/3/2006	2800	3.5	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		<2	4.5		<2	<2	<5
W-4	ug/L	11/9/2006	230	6.1	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		<2	5.1		<2	<2	<5
W-4	ug/L	2/8/2007	200	3.1	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		<2	4.7		<2	<2	<5
W-4	ug/L	5/10/2007	170	1.5	<2	<2	<2	<2	1.6	30	<5	<2	<2	<2		<2	3.8		<2	<2	1
W-4	ug/L	8/9/2007	280	1	<2	<2	<2	<2	2	18	<5	<2	<2	<2		<2	3.2		<2	<2	0.59
W-4	ug/L	11/7/2007	180	1.9	<0.36	<0.25	<0.6	<0.3	1.4	22	<0.41	<0.23	<0.26	<0.32		<0.27	3.6		0.36	<0.28	<0.3
W-4	ug/L	2/7/2008	210	4.4	<2	<2	<2	<2	<5	55	<1	<2	<2	<2		<1	4.4		<2	<2	<5
W-4	ug/L	2/7/2008	250	3.9	<2	<2	<2	<2	<5	50	<5	<2	<2	<2		<2	4		<2	<2	<5

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
W-4	ug/L	1/19/2009	140	0.51	<2	<2	<2	<2	<5	47	0.43	<2	<2	<2		<2	7.6		1	<2	1.8
W-4	ug/L	4/27/2009	92	<2	<2	<2	<2	<2	<5	34	<5	<2	<2	<2		<2	7.3		0.61	<2	1.9
W-4	UG/L	3/5/2010	600	1.5	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	3.7		<1.0	<0.50	7.4
W-4	UG/L	5/13/2010	700	4.3	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	3.1		<1.0	<0.50	5.4
W-4	UG/L	8/6/2010	570	68	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	4.0		<1.0	<0.50	7.2
W-4	UG/L	11/4/2010	980	180	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	4.8
W-4	UG/L	2/8/2011	1800	480	<0.50	1.2	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.0	<1.0	<1.0	<0.50	8.6
W-4	UG/L	4/14/2011	1400	460	0.59	1.2	<1.0	<0.50	1.1	38	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.2	<1.0	<1.0	<0.50	11
W-4	UG/L	8/25/2011	840	190	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	1.8
W-4	UG/L	11/14/2011	1200	390	<2.5	0.76	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-4	UG/L	2/6/2012	1100	410	<0.50	0.79	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	6.2
W-4	UG/L	5/7/2012	910	140	<0.50	<0.50	<1.0	<0.50	<1.0	21	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	4.1
W-4	UG/L	8/27/2012	910	<0.50	<0.50	<0.50	<1.0	<0.50	1.9	24	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	2.8
W-4	UG/L	11/5/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	6.3	2.5	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-4	UG/L	1/30/2013	160	28	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	1.2	<1.0	<1.0	<1.0	<1.0	1.5	<1.0	<1.0	<0.50	<1.0
W-4	UG/L	1/30/2013	190	43	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.1	<1.0	<1.0	<0.50	<1.0
W-4	UG/L	4/8/2013	360	18	<0.50	<0.50	<1.0	<0.50	2.8	77	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.3	<1.0	<1.0	<0.50	3.8
W-7	ug/L	8/4/2000	<500	<0.5	<1	<1			<1		<1	<1	<1	<1		<1	<0.5		1.2	<1	<0.5
W-7	ug/L	2/8/2001	<500	<0.5	<1	<1			<1		<10	<1	<1	<1		<1	<1		<1	<0.5	<0.5
W-7	ug/L	7/26/2001	<100	<0.5	<1	<1			<1		<10	<1	<1	<1		<1	<1		<1	<0.5	<0.5
W-7	ug/L	5/7/2002	<100	<0.5	<1	<1			<1	<10000	<10	<1	<1	<1		<1	<1		<1	<0.5	<0.5
W-7	ug/L	9/24/2002	<100	<0.5	<1	<1			<1	<10000	<10	<1	<1	<1		<1	<1		<1	<0.5	<0.5
W-7	ug/L	10/7/2005	<100	<0.5	<1	<1	<1	<1	<1	<10	<10	<1	<1	<1		<1	<1		<1	<0.5	<0.5
W-7	ug/L	2/16/2006	60.9	<1	<5	<5	<5	<5	<1	<10	<5	1.1	<5	<5		<5	<5		<5	<5	<5
W-7	ug/L	8/4/2006	<50	<2	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-7	ug/L	11/10/2006	<50	<2	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-7	ug/L	2/9/2007	<50	<2	<2	<2	2.6	<2	<5	<50	<5	2.2	<2	<2		<2	<2		<2	<2	<5
W-7	ug/L	5/8/2007	31	0.41	0.45	0.87	1.4	0.75	<5	<50	0.9	1.4	0.35	<2		<2	<2		0.41	<2	<5
W-7	ug/L	8/10/2007	<50	<2	<2	0.25	<2	<2	<5	<50	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-7	ug/L	11/6/2007	<30	<0.28	<0.36	<0.25	<0.6	<0.3	<0.32	<4.9	<0.41	<0.23	<0.26	<0.32		<0.27	<0.32		<0.27	<0.28	<0.3
W-7	ug/L	2/4/2008	<50	<2	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-7	ug/L	1/13/2009	<50	<2	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-7	ug/L	4/21/2009	<50	0.31	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		<2	<2		1.7	<2	<5
W-7	UG/L	3/4/2010	65	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		2.0	<0.50	<1.0
W-7	UG/L	5/17/2010	60	<0.50	<0.50	<0.50		0.51	<1.0	<10	2.3	<1.0	<1.0	<1.0		<1.0	<1.0		1.9	<0.50	<1.0
W-7	UG/L	8/4/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		2.6	<0.50	<1.0
W-7	UG/L	8/4/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		2.6	<0.50	<1.0
W-7	UG/L	11/3/2010	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.5	<0.50	<1.0
W-7	UG/L	2/2/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.4	<0.50	<1.0
W-7	UG/L	4/14/2011	<50	0.57	0.55	0.51	<1.0	0.57	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.8	<0.50	<1.0
W-7	UG/L	8/24/2011	<50	0.52	0.50	0.53	<1.0	0.53	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.8	<0.50	<1.0
W-7	UG/L	8/24/2011	<50	<0.50	<0.50	<0.50	<1.0	0.51	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.8	<0.50	<1.0
W-7	UG/L	11/10/2011	<50	<0.50	<0.50	0.56	<1.0	0.61	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.8	<0.50	<1.0
W-7	UG/L	2/8/2012	<50	<0.50	<0.50	0.57	<1.0	0.59	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.6	<0.50	<1.0
W-7	UG/L	5/9/2012	57	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.6	<0.50	<1.0
W-7	UG/L	8/29/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-7	UG/L	11/7/2012	<50	0.53	<0.50	0.64	<1.0	0.57	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
W-7	UG/L	2/1/2013	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-7	UG/L	4/4/2013	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-8	ug/L	8/4/2000	<500	2.8	<4.6	<1			<1		<1	<1	<1	<1		<1	<1		<1	<0.5	<0.5
W-8	ug/L	2/6/2001	NS	NS	NS	NS			NS		NS	NS	NS	NS		NS	NS		NS	NS	NS
W-8	ug/L	7/26/2001	180	0.67	<1	<1			<1		<1	<1	<1	<1		<1	<1		<1	<5	<0.5
W-8	ug/L	5/7/2002	180	0.51	<1	<1			<1	<10000	<10	<1	<1	<1		<1	<1		<1	<5	<0.5
W-8	ug/L	9/24/2002	<100	0.64	<1	<1			<1	<10000	<10	<1	<1	<1		<1	<1		<1	<5	<0.5
W-8	ug/L	7/1/2004	390	1.9J	1.8	0.72	0.92	<0.5	<5	<100	<5	<5	<5	<5		<5	<5		<5	<5	<5
W-8	ug/L	10/6/2005	220	0.52	<1	<1	<1	<1	<1	<10	<10	<1	<1	<1		<1	<1		<1	<0.5	<0.5
W-8	ug/L	2/16/2006	192	<1	<5	<5	<5	<5	<1	<10	<5	<5	<5	<5		<5	<5		<5	<5	<5
W-8	ug/L	8/4/2006	130	<2	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-8	ug/L	11/10/2006	210	<2	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-8	ug/L	2/9/2007	130	<2	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-8	ug/L	5/8/2007	110	0.49	0.73	0.33	<2	<2	<5	<50	<5	0.23	<2	<2		<2	<2		<2	<2	<5
W-8	ug/L	8/7/2007	170	0.49	0.82	0.44	<2	0.38	<5	<50	<5	0.3	<2	<2		<2	<2		<2	<2	<5
W-8	ug/L	11/6/2007	160	0.52	0.75	0.4	<0.6	0.3	<0.32	7.5	<0.41	<0.23	<0.26	<0.32		<0.27	<0.32		<0.27	<0.28	<0.3
W-8	ug/L	2/4/2008	160	0.46	0.81	0.39	<2	<2	<5	<50	<5	0.25	<2	<2		<2	<2		<2	<2	<5
W-8	ug/L	1/13/2009	120	<2	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-8	ug/L	4/21/2009	150	0.45	0.82	0.37	<2	<2	<5	<50	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-8	UG/L	3/4/2010	220	<0.50	0.85	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-8	UG/L	5/17/2010	200	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-8	UG/L	5/17/2010	210	<0.50	0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-8	UG/L	8/4/2010	110	<0.50	0.80	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	<1.0		<1.0	<0.50	<1.0
W-8	UG/L	11/4/2010	140	<0.50	0.60	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-8	UG/L	2/7/2011	130	<0.50	0.85	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-8	UG/L	4/21/2011	130	0.57	1.1	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-8	UG/L	4/21/2011	140	0.56	1.0	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-8	UG/L	9/1/2011	2000	0.57	0.77	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-8	UG/L	11/10/2011	110	<0.50	0.64	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-8	UG/L	2/7/2012	90	<0.50	0.73	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-8	UG/L	5/10/2012	180	<0.50	0.87	<0.50	<1.0	<0.50	<1.0	<10	2.9	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-8	UG/L	8/29/2012	190	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-8	UG/L	11/7/2012	62	0.50	0.75	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-8	UG/L	2/13/2013	<50	<0.50	0.87	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-8	UG/L	4/4/2013	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-9	ug/L	11/7/2006	<50	<2	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-9	ug/L	2/6/2007	67	<2	<2	<2	<2	<2	<5	<50	<5	<2	<2	<2		<2	<2		<2	<2	<5
W-9	ug/L	5/9/2007	50	<2	<2	<2	<2	<2	<5	17	<5	<2	<2	<2		<2	2		<2	<2	<5
W-9	ug/L	8/7/2007	38	<2	<2	<2	<2	<2	<5	22	<5	<2	<2	<2		0.31	3		<2	<2	<5
W-9	ug/L	11/6/2007	<30	<0.28	<0.36	<0.25	<0.6	<0.3	<0.32	19	<0.41	<0.23	<0.26	<0.32		0.31	3.8		<0.27	<0.28	<0.3
W-9	ug/L	2/5/2008	<50	<2	<2	<2	<2	<2	<5	23	0.5	<2	<2	<2		0.3	3.4		<2	<2	<5
W-9	ug/L	1/15/2009	46	<2	<2	<2	<2	<2	<5	18	<5	<2	<2	<2		<2	3.2		<2	<2	<5
W-9	ug/L	4/23/2009	36	<2	<2	<2	<2	<2	<5	18	<5	<2	<2	<2		<2	2.6		<2	<2	<5
W-9	UG/L	3/3/2010	<50	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	1.9		<1.0	<0.50	<1.0
W-9	UG/L	5/12/2010	80	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	2.8		<1.0	<0.50	<1.0
W-9	UG/L	8/4/2010	67	<0.50	<0.50	<0.50		<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0		<1.0	4.0		<1.0	<0.50	<1.0

Table III
Summary of Total Petroleum Hydrocarbon (TPH) and VOC Results
Former Powerine Refinery
Santa Fe Springs, CA
2Q2013

Location	Unit	Date	TPH-g	B	T	E	m/p-X	o-X	MTBE	TBA	NAP	1,2,4-TMB	1,3,5-TMB	PCE	TCE	t1,2-DCE	c1,2-DCE	1,1-DCE	1,1-DCA	1,2-DCA	VC
W-9	UG/L	11/3/2010	87	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	3.2	<1.0	<1.0	<0.50	<1.0
W-9	UG/L	2/2/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.5	<1.0	<1.0	<0.50	<1.0
W-9	UG/L	4/14/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	5.9	<1.0	<1.0	<0.50	<1.0
W-9	UG/L	8/24/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.4	<1.0	<1.0	<0.50	<1.0
W-9	UG/L	11/10/2011	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.1	<1.0	<1.0	<0.50	<1.0
W-9	UG/L	2/8/2012	59	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	13	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.8	<1.0	<1.0	<0.50	<1.0
W-9	UG/L	5/9/2012	89	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	29	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	2.3	<1.0	<1.0	<0.50	<1.0
W-9	UG/L	8/28/2012	70	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-9	UG/L	11/7/2012	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-9	UG/L	1/31/2013	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	<0.50	<1.0
W-9	UG/L	4/5/2013	<50	<0.50	<0.50	<0.50	<1.0	<0.50	<1.0	<10	<1.0	<1.0	<1.0	<1.0	<1.0	<1.0	1.1	<1.0	<1.0	<0.50	<1.0

NOTES:

PCE - Tetrachloroethylene

TCE - Trichloroethylene

c1,2-DCE - cis-1,2-Dichloroethene

t1,2-DCE - trans-1,2-Dichloroethene

1,1-DCE - 1,1-Dichloroethene

1,2-DCA - 1,2-Dichloroethane

1,3,5-TMB - 1,3,5-Trimethylbenzene

1,2,4-TMB - 1,2,4-Trimethylbenzene

VC - Vinyl Chloride

B- Benzene

T - Toluene

E - Ethylbenzene

X - Xylenes, total

nBUT - n-Butylbenzene

sBUT - sec-Butylbenzene

tBUT - tert-Butylbenzene

nPRO - n-Propylbenzene

1,1 DCA - 1,1-Dichloroethane

ISO-P - Isopropylbenzene

MC - Methylene Chloride

NAP - Naphthalene

TRIM - Trichlorofluoromethane

PMXY - p/m-Xylenes

OXYL -o-Xylene

DIPE - Diisopropyl Ether (DIPE)

MTBE - Methyl-tert-Butyl Ether (MTBE)

TBA - tert-Butyl Alcohol (TBA)

ND - Not Detected above laboratory detection limits

UG/L - Micrograms per litre

NA - Information not available

Table IV
Summary of Field Test Parameters
Former Powerine Refinery
Santa Fe Springs, California
2Q2013

Well ID	Sample Date	pH (SU)	DO (mg/L)	ORP (mV)
MW-104A	12/18/2009	7.31	5.31	3
MW-104A	3/3/2010	6.93	1.65	66
MW-104A	5/11/2010	8.06	NM	19
MW-104A	8/4/2010	7.65	2.32	205
MW-104A	11/3/2010	8.06	2.00	131
MW-104A	2/2/2011	8.46	3.05	136.4
MW-104A	4/14/2011	8.10	2.85	128.5
MW-104A	8/24/2011	7.53	4.47	19.6
MW-104A	11/10/2011	7.38	5.47	67
MW-104A	2/9/2012	8.79	2.42	-14.5
MW-104A	5/9/2012	8.18	4.36	-39.3
MW-104A	8/27/2012	7.69	1.96	51.9
MW-104A	11/6/2012	NM	NM	NM
MW-104A	1/28/2013	7.80	2.52	-43.6
MW-104A	4/5/2013	NM	NM	NM
MW-106A	12/17/2009	7.25	7.29	-112
MW-106A	3/5/2010	6.73	4.71	116
MW-106A	5/13/2010	8.06	7.90	-38
MW-106A	8/6/2010	8.05	4.52	210
MW-106A	11/4/2010	8.23	3.09	77
MW-106A	2/3/2011	NM	NM	NM
MW-106A	4/19/2011	NM	NM	NM
MW-106A	8/25/2011	7.67	2.98	-28.1
MW-106A	11/14/2011	7.03	4.74	33
MW-106A	2/3/2012	NM	NM	NM
MW-106A	8/24/2012	NM	NM	NM
MW-106A	11/6/2012	NM	NM	NM
MW-106A	1/28/2013	NM	NM	NM
MW-106A	4/4/2013	NM	NM	NM
MW-107A	12/17/2009	7.20	6.99	-276
MW-107A	3/5/2010	8.70	1.81	-307
MW-107A	5/13/2010	8.30	NM	-370
MW-107A	8/6/2010	8.10	3.25	-280
MW-107A	11/4/2010	8.16	2.04	-245
MW-107A	2/3/2011	8.49	3.42	-338
MW-107A	4/19/2011	8.02	1.93	-276.8
MW-107A	8/25/2011	7.82	2.68	-216.7
MW-107A	11/14/2011	7.19	3.73	-161.3
MW-107A	1/31/2012	8.88	2.6	-240

Table IV
Summary of Field Test Parameters
Former Powerine Refinery
Santa Fe Springs, California
2Q2013

Well ID	Sample Date	pH (SU)	DO (mg/L)	ORP (mV)
MW-107A	5/8/2012	8.40	2.34	-273.6
MW-107A	8/24/2012	8.12	2.89	-226.7
MW-107A	11/6/2012	8.27	2.38	-236.7
MW-107A	1/28/2013	7.96	2.25	-257.3
MW-107A	4/4/2013	8.25	2.25	-251.8
MW-503B	12/15/2009	6.92	7.78	-137
MW-503B	3/8/2010	7.33	3.38	-96
MW-503B	5/17/2010	8.18	1.79	-69
MW-503B	8/9/2010	7.60	2.72	147
MW-503B	11/8/2010	7.62	2.93	7
MW-503B	2/4/2011	7.96	2.16	-46
MW-503B	4/15/2011	7.61	1.74	-46.4
MW-503B	8/29/2011	7.50	2.57	-96.1
MW-503B	11/16/2011	6.76	3.01	-41.3
MW-503B	1/31/2012	8.50	3.06	-150.6
MW-503B	5/8/2012	7.73	2.46	-145.0
MW-503B	8/30/2012	8.05	2.50	-13.0
MW-503B	11/5/2012	8.00	2.06	96.5
MW-503B	1/30/2013	7.67	2.10	31.9
MW-503B	4/8/2013	7.72	2.46	-31.5
W-1	12/15/2009	7.62	7.10	-39
W-1	3/5/2010	7.51	3.15	-111
W-1	5/13/2010	8.07	2.02	-197
W-1	8/6/2010	7.52	3.22	-22
W-1	11/5/2010	8.13	2.75	38
W-1	2/4/2011	8.18	4.84	-63.7
W-1	4/14/2011	7.65	1.94	37.3
W-1	8/26/2011	7.47	3.16	-86
W-1	11/14/2011	7.08	2.9	-75.9
W-1	2/6/2012	7.99	2.87	-79.4
W-1	5/7/2012	7.85	3.03	-62.4
W-1	8/27/2012	7.90	2.69	-60.4
W-1	11/5/2012	7.82	2.47	-40.0
W-1	1/30/2013	7.64	3.07	66.5
W-1	4/3/2013	8.06	3.57	-96.5
W-4	12/15/2009	8.27	9.40	21
W-4	3/5/2010	7.09	3.41	-101
W-4	5/13/2010	8.00	3.87	-66
W-4	8/6/2010	7.74	3.48	16

Table IV
Summary of Field Test Parameters
Former Powerine Refinery
Santa Fe Springs, California
2Q2013

Well ID	Sample Date	pH (SU)	DO (mg/L)	ORP (mV)
W-4	11/4/2010	7.75	3.50	45
W-4	2/8/2011	7.67	5.53	-3.5
W-4	4/14/2011	7.79	4.47	107.8
W-4	8/25/2011	7.54	4.75	-92.5
W-4	11/14/2011	6.88	4.49	-47.3
W-4	2/6/2012	8.36	3.7	-53.2
W-4	5/7/2012	8.10	3.24	-54
W-4	8/27/2012	8.08	3.84	11.7
W-4	11/5/2012	8.18	3.95	20.2
W-4	1/30/2013	7.66	2.86	111.4
W-4	4/8/2013	7.91	3.97	-45.8
W-8	12/18/2009	10.11	7.07	-230
W-9	3/3/2010	7.53	5.66	69
W-9	5/12/2010	8.07	7.15	-175
W-9	8/4/2010	7.36	3.36	-60
W-9	4/5/2011	7.71	4.07	82.3
W-9	8/24/2011	7.62	4.9	-4.9
W-9	11/10/2011	NM	NM	NM
W-9	2/8/2012	8.32	3.95	61.8
W-9	5/9/2012	7.77	3.69	-49.5
W-9	8/28/2012	7.70	2.61	36.6
W-9	11/7/2012	NM	NM	NM
W-9	1/31/2013	7.49	2.37	13.1
W-9	4/5/2013	7.72	2.81	-93.6
W-10	12/18/2009	7.21	6.89	-97
W-10	3/8/2010	NM	NM	NM
W-10	5/17/2010	NM	NM	NM
W-10	8/9/2010	NM	NM	NM
W-10	11/3/2010	7.53	3.39	-10
W-10	11/8/2010	NM	NM	NM
W-10	2/2/2011	7.83	3.57	41.6
W-10	2/8/2011	7.28	5.51	-103
W-10	4/15/2011	NM	NM	NM
W-10	8/29/2011	7.14	2.7	-130.2
W-10	11/10/2011	NM	NM	NM
W-10	2/8/2012	NM	NM	NM
W-10	5/10/2012	NM	NM	NM
W-10	8/28/2012	NM	NM	NM
W-10	11/7/2012	NM	NM	NM

Table IV
Summary of Field Test Parameters
Former Powerine Refinery
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Well ID	Sample Date	pH (SU)	DO (mg/L)	ORP (mV)
W-10	1/28/2013	NM	NM	NM
W-10	4/1/2013	NM	NM	NM
W-11	12/8/2010	NM	NM	NM
W-11	2/4/2011	7.67	5.62	-119
W-11	4/15/2011	7.58	1.68	-77
W-11	8/29/2011	7.35	2.2	-125.7
W-11	11/14/2011	6.93	2.63	-148.6
W-11	2/8/2012	8.38	3.3	45.6
W-11	5/10/2012	7.84	2.75	-76.5
W-11	8/28/2012	7.50	1.56	-122.5
W-11	11/8/2012	7.92	1.75	24.7
W-11	1/31/2013	7.64	2.62	-120.3
W-11	4/5/2013	7.81	2.36	-69.2
W-12	12/18/2009	6.99	6.96	0
W-12	3/4/2010	7.53	3.15	-63
W-12	5/12/2010	7.87	NM	-180
W-12	8/5/2010	7.61	2.65	-100
W-12	11/4/2010	7.88	2.64	7
W-12	2/3/2011	8.28	2.85	-99
W-12	4/19/2011	7.77	2.10	15.2
W-12	8/25/2011	7.50	2.78	-58.5
W-12	11/14/2011	6.93	3.77	-34.7
W-12	2/8/2012	8.13	2.57	-113
W-12	5/9/2012	7.89	3.22	-74.5
W-12	8/30/2012	7.63	2.15	-98.7
W-12	11/8/2012	7.88	2.31	-42.6
W-12	1/31/2013	7.76	2.18	-70.3
W-12	4/2/2013	7.83	1.74	-98.7
W-14A	12/15/2009	7.65	7.76	-23.0
W-14A	3/1/2010	6.61	4.09	58.0
W-14A	5/10/2010	8.63	2.74	2.0
W-14A	8/2/2010	8.02	3.12	145.0
W-14A	11/1/2010	8.30	2.87	46.0
W-14A	1/31/2011	8.30	13.16	185.4
W-14A	4/4/2011	8.29	4.81	89.6
W-14A	8/22/2011	7.87	10.15	22.8
W-14A	11/7/2011	7.40	5.23	151.6
W-14A	1/30/2012	8.06	1.48	2.6
W-14A	8/20/2012	8.10	3.44	-76.9

Table IV
Summary of Field Test Parameters
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Well ID	Sample Date	pH (SU)	DO (mg/L)	ORP (mV)
W-14A	10/29/2012	8.23	3.01	22.5
W-14A	1/22/2013	8.18	5.24	226.9
W-14A	4/1/2013	8.35	2.98	-168.2
W-14B	12/15/2009	8.37	7.79	97
W-14B	3/1/2010	7.72	2.60	-5
W-14B	5/10/2010	8.43	3.00	-172
W-14B	8/2/2010	7.80	4.60	33
W-14B	11/1/2010	8.13	3.37	37
W-14B	1/31/2011	8.17	19.82	194
W-14B	4/4/2011	8.27	5.95	82.6
W-14B	8/22/2011	7.95	7.90	22.7
W-14B	11/7/2011	7.22	4.92	67.8
W-14B	1/30/2012	8.70	2.90	-133.7
W-14B	8/20/2012	8.27	4.00	-30.3
W-14B	10/29/2012	8.21	3.49	-18.2
W-14B	1/22/2013	7.96	4.24	130.4
W-14B	4/1/2013	8.28	2.90	-203.6
W-14C	12/15/2009	8.24	8.57	77.0
W-14C	3/1/2010	7.22	2.43	188.0
W-14C	5/10/2010	8.17	0.80	-77.0
W-14C	8/2/2010	7.60	3.55	128.0
W-14C	11/1/2010	7.89	3.15	49.0
W-14C	1/31/2011	7.88	10.85	188.0
W-14C	4/4/2011	7.98	3.27	51.3
W-14C	8/22/2011	7.76	4.24	-3.7
W-14C	11/7/2011	7.33	7.47	59.2
W-14C	1/30/2012	8.75	3.65	-65.2
W-14C	5/1/2012	8.18	4.07	41.5
W-14C	8/20/2012	8.18	4.95	5.1
W-14C	10/29/2012	8.16	3.77	-20.0
W-14C	1/22/2013	7.88	3.37	127.5
W-14C	4/1/2013	8.22	2.63	-181.5
W-15A	12/14/2009	7.31	9.15	85.0
W-15A	3/2/2010	7.12	2.67	202.0
W-15A	5/10/2010	7.90	NM	-228.0
W-15A	8/2/2010	7.39	1.96	-145.0
W-15A	11/1/2010	7.67	2.85	32.0
W-15A	2/1/2011	7.89	2.05	-33.0
W-15A	4/5/2011	8.00	2.60	-81.7

Table IV
Summary of Field Test Parameters
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Well ID	Sample Date	pH (SU)	DO (mg/L)	ORP (mV)
W-15A	8/23/2011	7.47	4.96	-148.7
W-15A	11/8/2011	(FPPH)	(FPPH)	(FPPH)
W-15A	2/2/2012	(FPPH)	(FPPH)	(FPPH)
W-15A	5/2/2012	8.06	3.26	-26.4
W-15A	8/21/2012	(FPPH)	(FPPH)	(FPPH)
W-15A	10/30/2012	(FPPH)	(FPPH)	(FPPH)
W-15A	1/23/2013	(FPPH)	(FPPH)	(FPPH)
W-15A	4/2/2013	(FPPH)	(FPPH)	(FPPH)
W-15B	12/14/2009	7.39	7.44	-58.0
W-15B	3/2/2010	7.61	2.39	94.0
W-15B	5/11/2010	8.09	4.36	-15.0
W-15B	8/3/2010	7.74	3.42	107.0
W-15B	11/2/2010	8.06	3.18	40.0
W-15B	2/1/2011	8.15	4.58	286.0
W-15B	4/5/2011	8.10	2.92	62.4
W-15B	8/23/2011	7.56	3.85	-2.1
W-15B	11/10/2011	7.10	3.07	28.3
W-15B	2/2/2012	8.17	2.31	-69.2
W-15B	5/2/2012	8.00	3.41	-11.0
W-15B	8/20/2012	8.10	5.08	64.6
W-15B	10/30/2012	8.21	2.80	123.6
W-15B	1/23/2013	7.75	2.74	135.0
W-15B	4/2/2013	8.16	1.88	-109.8
W-15C	12/14/2009	7.16	7.18	-53.0
W-15C	3/2/2010	7.33	2.27	148.0
W-15C	5/11/2010	8.16	4.73	-21.0
W-15C	8/3/2010	7.60	2.72	108.0
W-15C	11/2/2010	7.55	2.40	62.0
W-15C	2/1/2011	7.81	4.58	123.7
W-15C	4/5/2011	7.92	2.85	109.0
W-15C	8/23/2011	7.54	4.32	-2.4
W-15C	11/8/2011	7.32	6.00	119.4
W-15C	1/31/2012	8.72	3.11	-60.3
W-15C	5/2/2012	8.00	3.50	6.0
W-15C	8/21/2012	8.12	2.90	125.7
W-15C	10/30/2012	8.13	2.55	99.3
W-15C	1/23/2013	7.82	3.12	135.8
W-15C	4/2/2013	7.93	2.16	-107.1

Table IV
Summary of Field Test Parameters
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Well ID	Sample Date	pH (SU)	DO (mg/L)	ORP (mV)
W-16A	12/16/2009	7.62	6.90	-62.0
W-16A	3/5/2010	7.03	3.47	-5.0
W-16A	5/14/2010	8.28	2.23	-54.0
W-16A	8/9/2010	7.98	2.65	106.0
W-16A	11/5/2010	8.03	6.15	48.0
W-16A	2/7/2011	7.82	4.09	249.0
W-16A	4/18/2011	7.88	4.00	94.9
W-16A	8/26/2011	7.73	4.11	-73.4
W-16A	11/8/2011	7.07	4.36	77.6
W-16A	2/3/2012	8.49	3.67	-70.0
W-16A	5/3/2012	7.86	4.09	50.0
W-16A	8/22/2012	7.77	2.47	-77.5
W-16A	10/31/2012	8.15	4.03	113.1
W-16A	1/24/2013	7.77	3.30	64.6
W-16A	4/3/2013	7.80	2.83	-59.6
W-16B	12/16/2009	8.23	7.61	-184
W-16B	3/8/2010	8.15	3.20	-236
W-16B	5/14/2010	8.62	0.77	-310
W-16B	8/9/2010	8.01	2.88	-217
W-16B	11/5/2010	8.30	2.68	-119
W-16B	2/7/2011	8.12	3.54	-297
W-16B	4/18/2011	8.47	2.56	-247
W-16B	8/26/2011	8.01	2.72	-217.4
W-16B	11/8/2011	6.89	8.68	-63.8
W-16B	2/3/2012	9.21	2.55	-206.7
W-16B	5/3/2012	8.74	3.06	-194.3
W-16B	8/22/2012	8.62	2.90	-200.0
W-16B	10/31/2012	8.62	3.88	-189.5
W-16B	1/24/2013	7.96	2.53	-184.5
W-16B	4/3/2013	8.45	2.10	-198.3
W-16C	12/16/2009	8.15	7.12	-206
W-16C	3/8/2010	8.33	3.64	-237
W-16C	5/14/2010	8.68	NM	-295
W-16C	8/9/2010	8.02	2.57	-165
W-16C	11/5/2010	8.24	2.37	-72
W-16C	2/7/2011	8.03	4.34	-285
W-16C	4/18/2011	8.55	2.88	-249.5
W-16C	8/26/2011	7.81	2.71	-223.2
W-16C	11/9/2011	7.57	6.94	-185

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Well ID	Sample Date	pH (SU)	DO (mg/L)	ORP (mV)
W-16C	2/3/2012	8.84	2.51	-253.2
W-16C	5/3/2012	8.52	3.00	-205.8
W-16C	8/22/2012	8.30	2.60	-138.7
W-16C	10/31/2012	8.25	2.93	-185.2
W-16C	1/24/2013	8.05	3.20	-160.6
W-16C	4/3/2013	8.03	2.42	-169.2
W-17A	12/18/2009	8.02	7.10	30
W-17A	3/3/2010	6.67	5.41	74
W-17A	5/12/2010	8.25	0.88	-40
W-17A	8/4/2010	7.78	2.35	62
W-17A	11/3/2010	8.17	2.95	76
W-17A	2/2/2011	8.36	5.96	349
W-17A	4/20/2011	7.85	3.51	-5.8
W-17A	8/24/2011	7.85	3.23	2.6
W-17A	11/9/2011	7.19	4.78	-13
W-17A	2/7/2012	8.46	2.87	-20
W-17A	5/4/2012	8.20	3.45	-43.8
W-17A	8/23/2012	8.12	2.36	20.5
W-17A	11/1/2012	8.28	3.09	78.2
W-17A	1/25/2013	8.06	2.41	97.9
W-17A	4/9/2013	7.94	2.67	-27.8
W-17B	12/18/2009	8.49	7.18	-173
W-17B	3/3/2010	7.87	4.80	-197
W-17B	5/12/2010	8.35	NM	-313
W-17B	8/5/2010	7.96	2.31	-189
W-17B	11/3/2010	8.09	2.56	-25
W-17B	2/2/2011	8.43	3.45	-269
W-17B	4/20/2011	8.11	3.32	-168.5
W-17B	8/24/2011	7.88	3.41	-153.7
W-17B	11/9/2011	7.52	2.94	-136.4
W-17B	2/7/2012	8.65	2.50	-174.3
W-17B	5/4/2012	8.40	2.87	-118.7
W-17B	8/23/2012	8.25	2.13	-156.5
W-17B	11/1/2012	8.45	2.35	-97.2
W-17B	1/25/2013	8.14	2.81	-35.2
W-17B	4/9/2013	8.12	2.69	-155.6
W-17C	12/18/2009	8.79	8.74	-177
W-17C	3/4/2010	7.96	5.90	-209
W-17C	5/12/2010	8.49	3.03	-322

Table IV
Summary of Field Test Parameters
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Well ID	Sample Date	pH (SU)	DO (mg/L)	ORP (mV)
W-17C	8/5/2010	8.01	2.64	-167
W-17C	11/3/2010	8.16	2.79	-120
W-17C	2/2/2011	8.47	3.96	-301
W-17C	4/20/2011	8.26	2.08	-223.7
W-17C	8/24/2011	7.94	3.12	-201.7
W-17C	11/9/2011	7.43	3.36	-159.7
W-17C	2/7/2012	8.80	2.73	-226.4
W-17C	5/4/2012	8.50	2.56	-168.5
W-17C	8/23/2012	8.39	2.39	-177.5
W-17C	11/1/2012	8.48	2.87	-151.4
W-17C	1/25/2013	8.20	3.62	-166.8
W-17C	4/9/2013	8.25	2.36	-157.4
EW-1	2/3/2011	7.90	6.61	-258
EW-1	4/13/2011	8.15	2.86	-210
EW-1	8/29/2011	7.62	2.74	-293
EW-1	11/16/2011	(FPPH)	(FPPH)	(FPPH)
EW-1	2/6/2012	(FPPH)	(FPPH)	(FPPH)
EW-1	5/7/2012	(FPPH)	(FPPH)	(FPPH)
EW-1	8/24/2012	(FPPH)	(FPPH)	(FPPH)
EW-1	11/13/2012	(FPPH)	(FPPH)	(FPPH)
EW-1	1/29/2013	(FPPH)	(FPPH)	(FPPH)
EW-1	4/10/2013	(FPPH)	(FPPH)	(FPPH)
MW-701	2/4/2011	6.09	NM	NM
MW-701	4/11/2011	7.60	3.67	180.6
MW-701	8/30/2011	7.50	3.98	-31.2
MW-701	11/16/2011	6.90	2.93	25.9
MW-701	2/1/2012	8.18	4.3	-58.5
MW-701	5/11/2012	7.89	3.45	-8.8
MW-701	8/31/2012	7.97	4.00	28.7
MW-701	11/13/2012	7.88	3.00	161.0
MW-701	2/4/2013	7.84	4.20	120.5
MW-701	4/10/2013	7.84	3.55	36.6
MW-702	2/4/2011	6.04	NM	NM
MW-702	4/12/2011	7.70	3.29	103.1
MW-702	8/30/2011	7.34	3.23	-155.3
MW-702	11/16/2011	7.07	2.67	-172.7
MW-702	2/9/2012	7.89	4.73	-60.7
MW-702	5/11/2012	7.77	3.14	-99.9
MW-702	8/31/2012	7.76	3.48	-92.8

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Well ID	Sample Date	pH (SU)	DO (mg/L)	ORP (mV)
MW-702	11/13/2012	7.74	2.77	-116.3
MW-702	2/4/2013	7.60	3.34	-28.4
MW-702	4/10/2013	7.55	3.26	-26.9
MW-703	2/4/2011	6.25	NM	NM
MW-703	4/12/2011	7.57	3.53	132.4
MW-703	8/30/2011	7.30	4.2	-87.1
MW-703	11/17/2011	6.92	2.77	-98
MW-703	2/14/2012	8.11	4.07	-26.3
MW-703	5/11/2012	7.85	3.13	-72.6
MW-703	8/31/2012	7.68	3.20	-21.3
MW-703	11/14/2012	NM	NM	NM
MW-703	2/4/2013	7.75	3.50	122.6
MW-703	4/10/2013	7.87	3.75	-54.2
MW-704	2/9/2011	6.08	NM	NM
MW-704	4/13/2011	7.46	4.60	134.6
MW-704	8/31/2011	7.40	4.02	99.4
MW-704	11/17/2011	6.93	2.51	-148.8
MW-704	2/14/2012	7.80	4.2	-31.6
MW-704	5/14/2012	7.60	5.25	-30.0
MW-704	9/4/2012	7.87	2.85	31.7
MW-704	11/14/2012	NM	NM	NM
MW-704	2/5/2013	7.57	4.83	71.3
MW-704	4/15/2013	7.72	3.28	25.6
MW-705	2/4/2011	6.01	NM	NM
MW-705	4/12/2011	7.79	3.40	127.6
MW-705	8/31/2011	7.78	3.7	-55.5
MW-705	11/17/2011	7.04	3.16	-130.7
MW-705	2/14/2012	8.12	4.09	-57.6
MW-705	5/14/2012	7.88	2.50	-65.0
MW-705	9/4/2012	7.80	3.47	-28.4
MW-705	11/14/2012	NM	NM	NM
MW-705	2/5/2013	7.77	3.82	-46.8
MW-705	4/10/2013	7.73	2.78	21.8
MW-706	2/4/2011	6.21	NM	NM
MW-706	4/11/2011	7.99	4.02	158.7
MW-706	8/31/2011	7.76	3.03	-41.2
MW-706	11/18/2011	6.93	3.06	180.8
MW-706	2/14/2012	8.16	3.00	-52.7
MW-706	5/14/2012	7.87	2.77	-63.5

Table IV
Summary of Field Test Parameters
Former Powerine Refinery
Santa Fe Springs, California
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Well ID	Sample Date	pH (SU)	DO (mg/L)	ORP (mV)
MW-706	9/4/2012	7.84	3.24	18.2
MW-706	11/15/2012	8.04	3.31	-26.4
MW-706	2/5/2013	7.87	3.96	96.5
MW-706	4/15/2013	8.23	2.36	18.8
MW-707	2/4/2011	6.22	NM	NM
MW-707	4/8/2011	7.89	3.24	51.9
MW-707	9/1/2011	7.30	3.73	-9.4
MW-707	11/18/2011	6.89	2.8	11.3
MW-707	2/1/2012	8.19	3.1	-147
MW-707	5/15/2012	7.75	2.50	-72.6
MW-707	9/4/2012	7.55	3.26	-44.5
MW-707	11/15/2012	7.64	2.13	-88.8
MW-707	2/5/2013	7.62	3.58	13.1
MW-707	4/8/2013	7.67	3.23	-25.7
MW-708	2/4/2011	5.99	NM	NM
MW-708	4/6/2011	7.84	3.03	-119.8
MW-708	9/1/2011	7.51	3.45	-147.2
MW-708	11/18/2011	7.00	3.56	-161.3
MW-708	2/10/2012	8.09	2.75	-140.2
MW-708	5/15/2012	7.79	2.36	-136.1
MW-708	9/5/2012	7.78	2.39	-113.1
MW-708	11/16/2012	7.90	2.50	-133.6
MW-708	2/11/2013	7.62	3.47	-110.6
MW-708	4/11/2013	7.56	3.72	-28.5
MW-709	2/4/2011	6.27	NM	NM
MW-709	4/6/2011	8.08	3.74	149.6
MW-709	9/1/2011	7.38	2.97	-37
MW-709	11/21/2011	6.76	2.97	148.5
MW-709	2/10/2012	8.08	2.61	-57.1
MW-709	5/16/2012	7.70	3.12	9.3
MW-709	9/5/2012	7.82	2.07	-113.1
MW-709	11/16/2012	8.00	2.13	-78.2
MW-709	2/11/2013	7.61	3.00	59.4
MW-709	4/11/2013	7.76	2.62	88.1
MW-710	2/8/2011	6.18	NM	NM
MW-710	4/7/2011	7.88	3.54	97.7
MW-710	9/2/2011	6.87	3.68	-10.2
MW-710	11/21/2011	6.81	2.86	255.6
MW-710	2/1/2012	8.47	3.45	-64.8

Table IV
Summary of Field Test Parameters
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2Q2013

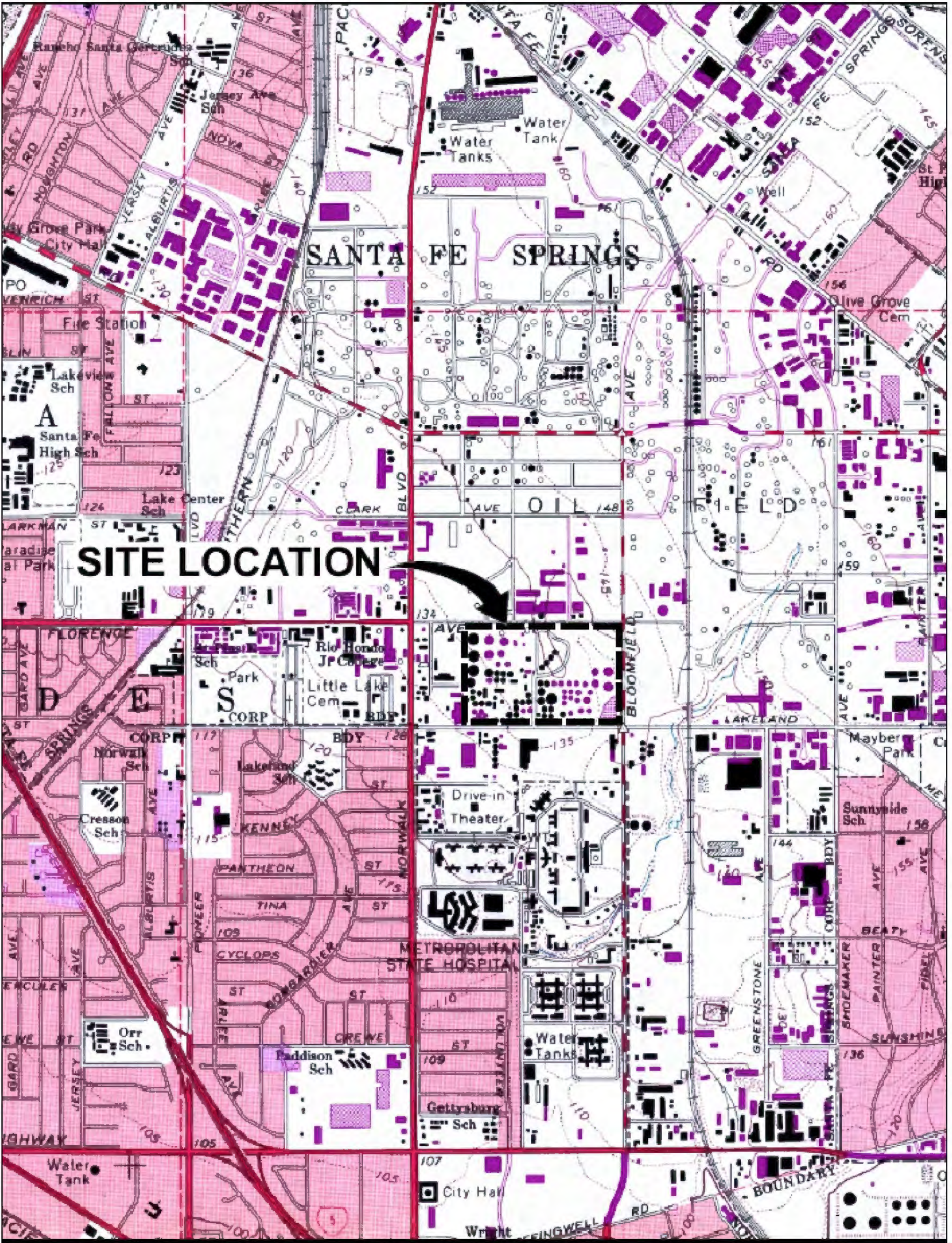
Well ID	Sample Date	pH (SU)	DO (mg/L)	ORP (mV)
MW-710	5/16/2012	7.80	4.04	21.5
MW-710	9/5/2012	7.85	2.32	30.5
MW-710	11/16/2012	7.97	3.57	43.4
MW-710	2/11/2013	7.63	3.13	94.0
MW-710	4/12/2013	7.77	2.66	62.3
MW-711	2/8/2011	5.99	NM	NM
MW-711	4/6/2011	7.91	3.39	-59.2
MW-711	9/2/2011	7.06	3.54	-99.8
MW-711	11/21/2011	6.87	2.95	-133.6
MW-711	2/10/2012	8.04	3.45	-96.7
MW-711	5/16/2012	7.73	2.37	-73.0
MW-711	9/5/2012	7.76	2.04	-175.4
MW-711	11/16/2012	7.77	2.66	-59.8
MW-711	2/11/2013	7.58	3.88	-66.4
MW-711	4/12/2013	7.67	2.32	-54.3
MW-712	2/7/2011	6.03	NM	NM
MW-712	4/7/2011	7.74	3.08	21.7
MW-712	9/2/2011	7.10	2.68	-59.7
MW-712	11/21/2011	6.90	2.65	-90.4
MW-712	2/13/2012	7.90	3.88	-83.5
MW-712	5/17/2012	7.71	2.80	-13.3
MW-712	9/6/2012	7.68	1.87	-42.0
MW-712	11/19/2012	7.83	2.26	-50.0
MW-712	2/12/2013	7.52	4.23	-5.3
MW-712	4/12/2013	7.59	2.21	-17.9
MW-713	2/7/2011	6.13	NM	NM
MW-713	4/8/2011	7.95	3.84	99.5
MW-713	9/2/2011	7.20	3.13	-51.4
MW-713	11/22/2011	6.98	3.07	-28.7
MW-713	2/13/2012	7.97	3.65	-77.7
MW-713	5/17/2012	7.70	3.11	-13.1
MW-713	9/6/2012	7.62	2.16	-120.7
MW-713	11/19/2012	7.79	2.72	-139.5
MW-713	2/12/2013	7.52	3.73	-101.8
MW-713	4/11/2013	7.66	2.95	-122.9
MW-714	2/8/2011	6.20	NM	NM
MW-714	4/7/2011	7.92	3.53	33.6
MW-714	9/2/2011	7.21	3.15	-63.4
MW-714	11/22/2011	6.96	2.77	-24.2

Table IV
Summary of Field Test Parameters
Former Powerine Refinery
Santa Fe Springs, California
2Q2013

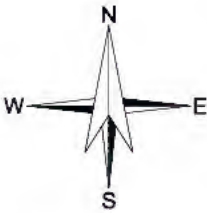
Well ID	Sample Date	pH (SU)	DO (mg/L)	ORP (mV)
MW-714	2/13/2012	8.05	4.32	-70.5
MW-714	5/17/2012	4.60	3.00	-10.7
MW-714	9/6/2012	7.66	2.58	-50.0
MW-714	11/19/2012	7.81	3.04	-98.7
MW-714	2/12/2013	7.58	4.77	-24.7
MW-714	4/11/2013	7.75	3.05	-54.3
MW-715	2/14/2011	7.50	NM	NM
MW-715	4/8/2011	7.78	2.59	16.3
MW-715	9/2/2011	7.15	3.2	-89.8
MW-715	11/22/2011	6.90	2.73	-125.4
MW-715	2/1/2012	8.32	2.87	-174.2
MW-715	5/17/2012	4.20	2.58	-50.5
MW-715	9/6/2012	7.66	1.97	-98.9
MW-715	11/19/2012	7.85	3.62	-134.5
MW-715	2/12/2013	7.65	3.75	-135.3
MW-715	4/12/2013	7.81	2.58	-124.9

Notes:

DO dissolved oxygen
mg/L milligram(s) per liter
mV millivolts
ORP oxidation-reduction potential
SU standard units
NM Not Measured



SOURCE OF BASE MAP
U.S. GEOLOGICAL SURVEY, 7.5 MIN QUAD., WHITTIER, CA. 1965, PHOTOREVISED 1981



SCALE: NOT TO SCALE

FORMER POWERINE REFINERY
12345 LAKELAND ROAD
SANTA FE SPRINGS, CALIFORNIA

SITE LOCATION MAP



FIGURE
1

FX-9 Wells

FX Wells

FX-9 Wells

FX-9 Wells

FX Wells

FX-9 Wells

FX Wells

FX-9 Wells

FX-9 Wells

FX-9 Wells

FX-9 Wells

FX-9 Wells

FX-9 Wells

Appendix A

Table 2- Groundwater Level Measurements
FS & AW

Well ID	Date 3.29.2013 2Q2013	Total Depth (feet BTOC)	Depth to Groundwater (feet BTOC)	Depth To FPPH FPPH (feet BTOC)	FPPH Thickness (feet)	Top of Casing Elevation (feet msl)	Groundwater Elevation (feet msl)	1Q2012
EW-1	FP	113	106.62	105.74	0	144.78		105.72 / 107.52
MW-101		90.72	—	NA	0	135.23		DRY
MW-103		94.7	—	NA	0	136.95		DRY
MW-104A	3-29-13	100.08	92.84	NA	0	143.39		93.42
MW-105		100.47	—	NA	0	138.63		DRY
MW-106A		110	104.24	NA	0	152.51		105.78
MW-107A		109.49	104.03	NA	0	146.71		104.67
MW-201		101.6	—	NA	0	132.91		DRY
MW-202		92.55	—	NA	0	137.89		DRY
MW-203		102.3	—	NA	0	143.43		DRY
MW-204		103.1	—	NA	0	142.18		DRY
MW-205		98.27	—	NA	0	138.04		DRY
MW-501A		93.27	—	NA	0	128.7		DRY
MW-502		100.59	—	NA	0	128.3		DRY
MW-503B		108.67	101.60		0	129.96		100.31
MW-504		95.76	—	NA	0	134.51		DRY
MW-600A		92.7	—	NA	0	120.34		DRY
MW-601A		89.9	—	NA	0	126.53		DRY
MW-603		97.6	—	NA	0	118.54		DRY
MW-604		103.2	—	NA	0	138.16		DRY
MW-605		93.98	—	NA	0	114.54		DRY
MW-606		99.05	—	NA	0	113.89		DRY
MW-607		107.05	—	NA	0	126.03		DRY
W-1		129.61	109.21	NA	0	142.89		109.91
W-10		110.21	97.98	NA	0	139.99		97.63
W-11		112.61	98.77		0	141.29		98.90
W-12		116.1	103.66	NA	0	144.42		103.39
W-14 A		112	99.78	NA	0	114.71		92.37
W-14 B		167	93.69	NA	0	114.79		90.35
W-14 C		195	13.91	NA	0	114.8		90.60
W-15 A	FP	125.7	113.59	111.08 NA	0	127.6		110.40
W-15 B		155.6	111.53	NA	0	127.62		110.11
W-15 C		197.34	111.85	NA	0	127.62		109.77
W-16 A		123.12	112.57	NA	0	147.61		113.40
W-16 B		160.25	117.92	NA	0	147.68		109.46
W-16 C		196.3	117.70	NA	0	147.67		109.12
W-17 A		108.3	97.25	NA	0	141.37		96.96
W-17 B		169.6	104.01	NA	0	141.34		98.15
W-17 C		200	104.02	NA	0	141.34		98.19
W-3A		111.73	—	NA		124		DRY
W-4		129.71	110.43	NA	0	142.38		111.13
W-7		NM	93.10	NA	0	NM		83.12
W-8		NM	77.20	NA	0	NM		67.75
W-9		110.37	90.93	NA	0	139.12		92.58
MW-701		130	99.91	NA	0			98.85
MW-702		130	99.52	NA	0			98.74
MW-703		130	101.17	NA	0			100.23
MW-704		130	102.90	NA	0			102.11
MW-705		130	103.90	NA	0			103.39
MW-706		130	100.48	NA	0			100.00
MW-707		130	98.46	NA	0			96.96
MW-708	FP	130	97.82	97.55	0			96.46 (no FP)
MW-709		130	107.65	NA	0			109.88
MW-710		130	96.25	NA	0			93.67
MW-711		130	103.00	NA	0			101.00
MW-712		130	97.95	NA	0			98.70
MW-713		130	105.59	NA	0			104.90
MW-714		142	106.18	NA	0			104.52
MW-715		134	97.98	NA	0			96.06

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013** *4-10-13*

WELL NO. **EW-1** Walker
 SAMPLED BY: **Frane Sasic**

WELL NOTES: **FPPH + strong vapors**

WELL CONDITION: *OK*

WEATHER CONDITIONS: *Clear/sunny/light breeze (~72°F)*

PURGING AND SAMPLING EQUIPMENT:
 YSI 556
 Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL <i>113.00</i>	(ft.)
DEPTH TO WATER	(ft.)
HEIGHT OF WATER COLUMN	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. =	(gal)
PURGE VOLUME x 3 =	(gal)
PRODUCT THICKNESS <i>106.62 (DW) - 105.74 (DTFP) =</i>	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. (s/cm)	Turbidity NTUs	DO mg/L	Temperature (F / C)	TDS	ORP	Color	Odor
/	/	/	<i>* FPPH *</i>	/	-	/	/	-	/	/	/
					-			-			
					-			-			

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	<i>4-10-13</i>	ice	8260B - VOCs + Oxys	VOAs	3	HCL	<i>Top-down FPPH skin</i> <i>No GW parameter measurements taken due to free product.</i> <i>50 gals purged total, of which ~13 gal is nearly black colored FP of low viscosity.</i>
1	<i>1000</i>	ice	8015M - TPH-g	VOAs	3	HCL	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(ft) \times 7.48 \text{ gal./ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

Murex Environmental Inc.

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GROUNDWATER SAMPLING LOG

PAGE 1 OF 2

PROJECT NAME: **CENCO**
PROJECT NO.: **1003-001-300**
DATE: **2Q2013 4-3-13**

WELL NO. **W-1** Walker
SAMPLED BY: **Frane Sasic**

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER	4" (inches)
DEPTH OF WELL	129.61 (ft.)
DEPTH TO WATER	109.21 (ft.)
HEIGHT OF WATER COLUMN	20.40 (ft.)
CASING VOLUME*	Hgt. x 0.163 Gal./Ft. = 13.464 (gal)
PURGE VOLUME	0.66 x 3 = 40.392 (gal)
PRODUCT THICKNESS	(ft.)

WELL NOTES:
WELL CONDITION:

OK

WEATHER CONDITIONS:

Mostly sunny w/ light breeze
(~80°F)

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. uS/cm	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
1456	5	VAC TRUCK	8.12	2.601	-	3.73	24.21	1.690	-99.3	Clear	Strong
1459	10		8.10	2.609	-	3.48	24.47	1.696	-101.5	Clear	Strong
1502	15	↓	8.13	2.613	-	3.69	24.24	1.697	-100.6	Clear	Strong

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	4-3-13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	Tip for air-line hose fitting is missing so we cannot use air assist to purge. Will try to purge w/o it. LL-W1-040313 @ 15:48
1	1548	ice	8015M - TPH-g	VOAs	3	HCL	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

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Page 2 of 2

WELL NO. *W-1*

SAMPLED BY: Frane Sosic

DATE: 4-3-13 (2Q2013)

[illegible]

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013 4-8-13**

WELL NO. **W-4** Walker
 SAMPLED BY: **Frane Sosic**

WELL NOTES:

WELL CONDITION:

OK

WEATHER CONDITIONS:

Cloudy and cool (~63°F) AM
Clear/Sunny / gusty winds (~73°F) PM

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL 129.71	(ft.)
DEPTH TO WATER 110.43	(ft.)
HEIGHT OF WATER COLUMN 19.28	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = 12.7248	(gal)
PURGE VOLUME 0.66 x 3 = 38.1744	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. $\mu S/cm$	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS μL	ORP μV	Color	Odor
9:29	5	VACTRUCK	8.08	2.307	--	2.76	21.91	1.498	-59.6	Green	Mild
9:35	10		8.12	2.280	--	2.78	21.02	1.482	-66.9	Green	Mild
9:49	15		7.95	2.262	--	3.01	20.25	1.470	-61.1	Light grey	Mild

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	4-8-13	ice	8260B - VOCs + OxyS	VOAs	3	HCL	Purge slowed down significantly after 10 gal.
1	1200	ice	8015M - TPH-g	VOAs	3	HCL	W-4 went dry ~ 23 gals. / allowed to re-charge prior to sampling:
							LL-W4-040813 @ 12:00
							* Stinger needed repairs *

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(ft) \times 7.48 \text{ gal./ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

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Page 2 of 2

PROJECT NO.: 1003-001-300

DATE: 4-8-13 (202013)

WELL NO.

SAMPLED BY: Frane Sosic

[illegible]

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013** *4-4-13*

WELL NO. **W-7** Lakeland
 SAMPLED BY: **Frane Sosic**

WELL NOTES: No purge well (sample in any order)

WELL CONDITION: GOOD

WEATHER CONDITIONS: Hazy + humid (~75°F)

PURGING AND SAMPLING EQUIPMENT:
 YSI 556
 Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER <i>~ 18"</i>	(inches)
DEPTH OF WELL	(ft.)
DEPTH TO WATER	(ft.)
HEIGHT OF WATER COLUMN	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. =	(gal)
PURGE VOLUME x 3 =	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. (s/cm)	Turbidity NTUs	DO mg/L	Temperature (F / C)	TDS	ORP	Color	Odor

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
<i>1</i>	<i>4-3-13</i>	<i>ice</i>	<i>8260B - VOCs + Oxys</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	<i>LL-W7-040413 @ 11:13</i>
<i>1</i>	<i>11:13</i>	<i>ice</i>	<i>8015M - TPH-g</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(ft) \times 7.48 \text{ gal./ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

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GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013 4-4-13**

WELL NO. **W-8** Lakeland
 SAMPLED BY: **Frane Sasic**

WELL NOTES: No purge well (sample in any order)

WELL CONDITION: GOOD

WEATHER CONDITIONS: Hazy + humid (~77°F)

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER ~18"	(inches)
DEPTH OF WELL	(ft.)
DEPTH TO WATER	(ft.)
HEIGHT OF WATER COLUMN	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. =	(gal)
PURGE VOLUME x 3 =	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. (s/cm)	Turbidity NTUs	DO mg/L	Temperature (F / C)	TDS	ORP	Color	Odor

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	4-4-13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	LL - W8-040413 @ 11:43
1	1143	ice	8015M - TPH-g	VOAs	3	HCL	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(ft) \times 7.48 \text{ gal./ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

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GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013 4-5-13**

WELL NO. **W-9** Lakeland
 SAMPLED BY: **Frane Sasic**

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 2"	(inches)
DEPTH OF WELL 110.00	(ft.)
DEPTH TO WATER 90.93	(ft.)
HEIGHT OF WATER COLUMN 19.07	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = 3.10841	(gal)
PURGE VOLUME x 3 = 9.32523	(gal)
PRODUCT THICKNESS	(ft.)

WELL CONDITION:
NOT GOOD

WEATHER CONDITIONS:
Scattered clouds w/ slight wind (~68°F)

PURGING AND SAMPLING EQUIPMENT:
 YSI 556
 Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. $\mu\text{S/cm}$	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS μS	ORP mV	Color	Odor
926	3	VAC TRUCK	7.99	2.772	—	3.94	22.01	1.802	-76.9	Rusty Flakes	Ugh
936	6	I	7.68	2.797	—	2.47	21.96	1.817	-78.7	Cloudy	Slight
949	9	I	7.72	2.789	—	2.81	20.93	1.810	-93.6	Cloudy	Slight

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	4-5-13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	LL - W9 - 04/05/13 @ 10:12
1	12:12	ice	8015M - TPH-g	VOAs	3	HCL	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal./ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013** *4-1-2013*

WELL NO. **W-10** Lakeland

SAMPLED BY: **Frane Sosis**

WELL NOTES: Slow recharge: purge 1 day prior to sample collection

WELL CONDITION: OK

WEATHER CONDITIONS:

Mostly sunny w/ light breeze (~75°F)

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 2"	(inches)
DEPTH OF WELL <i>110.00</i>	(ft.)
DEPTH TO WATER <i>97.98</i>	(ft.)
HEIGHT OF WATER COLUMN <i>12.02</i>	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = <i>1.95926</i>	(gal)
PURGE VOLUME x 3 = <i>5.87718</i>	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. (s/cm)	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS	ORP	Color	Odor
/	/	/	<i>* VOLUME *</i>	/	/	/	/	/	/	/	/

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
<i>1</i>	<i>4-1-13</i>	<i>ice</i>	<i>8260B - VOCs + Oxys</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	<i>*NO PARAMETERS TAKEN DUE TO LACK OF GW VOLUME FOR READING* W-10 went dry ~ 2 gallons. LL-W10-040113 @</i>
<i>1</i>		<i>ice</i>	<i>8015M - TPH-g</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal./ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

Murex Environmental Inc.

2640 Walnut Ave, Unit F, Tustin, CA 92780 | 714.508.0800 ph | 714.508.0880 fx | www.murexenv.com

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013 4-5-13**

WELL NO. **W-11** Lakeland

SAMPLED BY: **Frane Sosis**

WELL NOTES: **Historically contained FPPH**

WELL CONDITION:

GOOD

WEATHER CONDITIONS:

**Mostly sunny w/ some clouds + light wind
(~74°F)**

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 2"	(inches)
DEPTH OF WELL 113.00	(ft.)
DEPTH TO WATER 98.97	(ft.)
HEIGHT OF WATER COLUMN 14.03	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = 2.28689	(gal)
PURGE VOLUME x 3 = 6.86067	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. µS/cm	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
1357	5	VAC TRUCK	7.77	1.991	—	2.01	23.85	1.290	30 to -52.6	light grey	Strong
1411	10	↓	7.84	1.978	—	1.90	23.15	1.287	-65.6	cloudy	Strong
1433	15	↓	7.81	1.991	—	2.36	22.78	1.294	-69.2	Clear	Strong

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	4-5-13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	LL-W11-040513 @ 15:00
1	15:00	ice	8015M - TPH-g	VOAs	3	HCL	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal./ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

Murex Environmental Inc.

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GROUNDWATER SAMPLING LOG

PROJECT NAME: CENCO
 PROJECT NO.: 1003-001-300
 DATE: 2Q2013 4-2-2013

WELL NO. W-12 Lakeland
 SAMPLED BY: Frane Susic

WELL NOTES: May 2013

WELL CONDITION: GOOD

WEATHER CONDITIONS: Mostly sunny w/ light breeze
(~76°F)

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER <u>2"</u>	(inches)
DEPTH OF WELL <u>116.00</u>	(ft.)
DEPTH TO WATER <u>103.66</u>	(ft.)
HEIGHT OF WATER COLUMN <u>12.34</u>	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = <u>2.01142</u>	(gal)
PURGE VOLUME x 3 = <u>6.03426</u>	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. μS/cm	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
<u>1510</u>	<u>5</u>	<u>VAC TRUCK</u>	<u>7.77</u>	<u>2.117</u>	<u>/</u>	<u>2.46</u>	<u>24.52</u>	<u>1.373</u>	<u>-72.3</u>	<u>Light grey</u>	<u>Strong</u>
<u>1518</u>	<u>10</u>	<u> </u>	<u>7.81</u>	<u>2.159</u>	<u>/</u>	<u>1.83</u>	<u>24.41</u>	<u>1.402</u>	<u>-97.4</u>	<u>Cloudy</u>	<u>Mild</u>
<u>1530</u>	<u>15</u>	<u> </u>	<u>7.83</u>	<u>2.122</u>	<u>/</u>	<u>1.74</u>	<u>24.63</u>	<u>1.380</u>	<u>-98.7</u>	<u>Cloudy</u>	<u>Mild</u>

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
<u>1</u>	<u>4-2-13</u>	<u>ice</u>	<u>8260B - VOCs + Oxys</u>	<u>VOAs</u>	<u>3</u>	<u>HCL</u>	<u>LL - W12 - 040213 @ 1600</u>
<u>1</u>	<u>1600</u>	<u>ice</u>	<u>8015M - TPH-g</u>	<u>VOAs</u>	<u>3</u>	<u>HCL</u>	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013** *4-1-2013*

WELL NO. **MW-14A** Hospital
 SAMPLED BY: **Frane Sasic**

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 2"	(inches)
DEPTH OF WELL <i>112.00</i>	(ft.)
DEPTH TO WATER <i>94.78</i>	(ft.)
HEIGHT OF WATER COLUMN <i>17.22</i>	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = <i>2.80686</i>	(gal)
PURGE VOLUME x 3 = <i>8.42058</i>	(gal)
PRODUCT THICKNESS	(ft.)

WELL CONDITION:

OK

WEATHER CONDITIONS:

Overcast (~60°F)

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. <i>µS/cm</i>	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
<i>1000</i>	<i>5</i>	<i>VACUUM</i>	<i>8.27</i>	<i>1.605</i>	<i>/</i>	<i>3.37</i>	<i>22.3</i>	<i>1.047</i>	<i>-156.3</i>	<i>Cloudy</i>	<i>None</i>
<i>1037</i>	<i>10</i>	<i>TRUCK</i>	<i>8.35</i>	<i>1.623</i>	<i>/</i>	<i>2.98</i>	<i>21.91</i>	<i>1.055</i>	<i>-168.2</i>	<i>Clear</i>	<i>None</i>

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
<i>1</i>	<i>4-1-13</i>	<i>ice</i>	<i>8260B - VOCs + Oxys</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	<i>LL-14A 04/01/13 @ 10:45</i>
<i>1</i>	<i>@</i>	<i>ice</i>	<i>8015M - TPH-g</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	
<i>1</i>	<i>10:45</i>	<i>ice</i>	<i>8015 - Methane</i>	<i>VOAs</i>	<i>3</i>	<i>NONE</i>	
<i>1</i>	<i>-11-</i>	<i>ice</i>	<i>Total Alkalinity</i>	<i>250 ml poly</i>	<i>1</i>	<i>NONE</i>	
<i>1</i>	<i>-11-</i>	<i>ice</i>	<i>300 IC: nitrate, sulfate</i>	<i>250 ml poly</i>	<i>1</i>	<i>NONE</i>	
<i>1</i>	<i>-11-</i>	<i>ice</i>	<i>Ferrous Iron</i>	<i>250 ml poly</i>	<i>1</i>	<i>HNO₃</i>	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013** *4-1-2013*

WELL NO. **MW-14B** Hospital
 SAMPLED BY: **Frane Sasic**

WELL INFORMATION		
TOP OF CASING ELEV.		(ft.)
WELL DIAMETER	2"	(inches)
DEPTH OF WELL	<i>165.00</i>	(ft.)
DEPTH TO WATER	<i>93.69</i>	(ft.)
HEIGHT OF WATER COLUMN	<i>71.31</i>	(ft.)
CASING VOLUME*	Hgt. x 0.163 Gal./Ft. = <i>11.62353</i>	(gal)
PURGE VOLUME	x 3 = <i>34.87059</i>	(gal)
PRODUCT THICKNESS		(ft.)

WELL CONDITION:

OK

WEATHER CONDITIONS:

Scattered clouds / light breeze (20°F)

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. µS/cm	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
<i>1114</i>	<i>5</i>	<i>VACUUM</i>	<i>8.42</i>	<i>1.638</i>	<i>—</i>	<i>3.81</i>	<i>19.97</i>	<i>1.064</i>	<i>-199.0</i>	<i>Clear</i>	<i>Slight</i>
<i>1142</i>	<i>10</i>	<i>TIRUCK</i>	<i>8.31</i>	<i>1.624</i>	<i>—</i>	<i>2.77</i>	<i>20.36</i>	<i>1.056</i>	<i>-201.3</i>	<i>Clear</i>	<i>Slight</i>
<i>1210</i>	<i>15</i>	<i>↓</i>	<i>8.30</i>	<i>1.616</i>	<i>—</i>	<i>3.14</i>	<i>19.82</i>	<i>1.050</i>	<i>-200.7</i>	<i>Clear</i>	<i>Slight</i>
<i>1240</i>	<i>20</i>	<i>↓</i>	<i>8.28</i>	<i>1.605</i>	<i>—</i>	<i>2.90</i>	<i>21.11</i>	<i>1.043</i>	<i>-203.6</i>	<i>Clear</i>	<i>Slight</i>
Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES: <i>Very slow purging well</i>				
<i>1</i>	<i>4-1-13</i>	<i>ice</i>	<i>8260B - VOCs + Oxys</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	<i>LL-14B-040113 @ 12:48</i>				
<i>1</i>	<i>1248</i>	<i>ice</i>	<i>8015M - TPH-g</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>					

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(ft) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

GROUNDWATER SAMPLING LOG

PAGE 1 of 2

PROJECT NAME: **CENCO**
PROJECT NO.: **1003-001-300**
DATE: **2Q2013** *4-1-2013*

WELL NO. **MW-14C** Hospital
SAMPLED BY: **Frane Sasic**

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 2"	(inches)
DEPTH OF WELL <i>195.00</i>	(ft.)
DEPTH TO WATER <i>93.91</i>	(ft.)
HEIGHT OF WATER COLUMN <i>101.09</i>	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = <i>16.47767</i>	(gal)
PURGE VOLUME x 3 = <i>49.43301</i>	(gal)
PRODUCT THICKNESS	(ft.)

WELL CONDITION:

OK

WEATHER CONDITIONS:

Scattered clouds / light breeze (~75°F)

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp. Cond. $\mu\text{S/cm}$	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS <i>g/L</i>	ORP	Color	Odor
<i>1256</i>	<i>5</i>	<i>VACUUM</i>	<i>8.36</i>	<i>1.633</i>	<i>/</i>	<i>7.62</i>	<i>25.00</i>	<i>1.113</i>	<i>-211.4</i>	<i>Cloudy</i>	<i>Slight</i>
<i>1320</i>	<i>10</i>	<i>TRUCK</i>	<i>8.25</i>	<i>1.620</i>	<i>/</i>	<i>5.41</i>	<i>24.82</i>	<i>1.086</i>	<i>-196.5</i>	<i>Cloudy</i>	<i>Slight</i>
<i>1337</i>	<i>15</i>	<i>↓</i>	<i>8.21</i>	<i>1.612</i>	<i>/</i>	<i>3.60</i>	<i>24.15</i>	<i>1.067</i>	<i>-189.2</i>	<i>Cloudy</i>	<i>Slight</i>

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
<i>1</i>	<i>4-1-13</i>	<i>ice</i>	<i>8260B - VOCs + Oxys</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	<i>LL-14C-040113 @ 15:56</i>
<i>1</i>		<i>ice</i>	<i>8015M - TPH-g</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

Page 2 of 2

PROJECT NO.: 1003-001-300

DATE: 4-1-2013 (202013)

WELL NO. 14-C

SAMPLED BY: Frane Sosic

[illegible]

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013 4-2-2013**

WELL NO. **MW-15A** Hospital
 SAMPLED BY: **Frane Sasic**

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 2"	(inches)
DEPTH OF WELL 125.00	(ft.)
DEPTH TO WATER FPPH 111.08	(ft.)
HEIGHT OF WATER COLUMN 13.92	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = 2.26896	(gal)
PURGE VOLUME x 3 = 6.80	(gal)
PRODUCT THICKNESS 112.59(DTW) - 111.08(FP) = 2.51 ON 3/29/13	ft.)

Well Notes: **Contains feet of FPPH**

WELL CONDITION: **OK**

WEATHER CONDITIONS: **Cloudy w/ light wind (~60°F)**

PURGING AND SAMPLING EQUIPMENT:
 YSI SS6
 Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. (s/cm)	Turbidity NTUs	DO mg/L	Temperature (F / C)	TDS	ORP	Color	Odor
/	/	/	* FPPH *	/	/	/	/	/	/	/	/

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES: * NO PARAMETERS TAKEN DUE TO FREE PRODUCT *
1	4:2:13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	FPPH skin until we achieve steady GW flow: 15' of stinger removed initially, then stinger is lowered slowly to fluid level. Approx. 60 gal purged total: ~25 pure translucent brown FPPH, ~5 gal. of emulsified layer and ~30 gal. of GW. LL-15A-040213 @ 11:29
1	@	ice	8015M - TPH-g	VOAs	3	HCL	
1	11:29	ice	8015 - Methane	VOAs	3	NONE	
1	-11-	ice	Total Alkalinity	250 ml poly	1	NONE	
1	-11-	ice	300 IC: nitrate, sulfate	250 ml poly	1	NONE	
1	-11-	ice	Ferrous Iron	250 ml poly	1	HNO ₃	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(ft) \times 7.48 \text{ gal./ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013** *4-2-2013*

WELL NO. **MW-15B** Hospital
 SAMPLED BY: **Frane Sosic**

WELL INFORMATION		
TOP OF CASING ELEV.		(ft.)
WELL DIAMETER	2"	(inches)
DEPTH OF WELL	<i>156.00</i>	(ft.)
DEPTH TO WATER	<i>111.53</i>	(ft.)
HEIGHT OF WATER COLUMN	<i>44.47</i>	(ft.)
CASING VOLUME*	Hgt. x 0.163 Gal./Ft. = <i>7.24861</i>	(gal)
PURGE VOLUME	x 3 = <i>21.74583</i>	(gal)
PRODUCT THICKNESS		(ft.)

WELL CONDITION:

OK

WEATHER CONDITIONS:

*Mostly sunny w/light breeze
(~73°F)*

PURGING AND SAMPLING EQUIPMENT:

YSI S56

Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. <i>µmS/cm</i>	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS <i>g/L</i>	ORP <i>mV</i>	Color	Odor
<i>1203</i>	<i>5</i>	<i>VACUUM</i>	<i>8.40</i>	<i>2.099</i>	—	<i>2.38</i>	<i>21.61</i>	<i>1.366</i>	<i>-89.6</i>	<i>Light grey</i>	<i>Strong</i>
<i>1215</i>	<i>10</i>	<i>TRUCK</i>	<i>8.26</i>	<i>2.070</i>	—	<i>1.90</i>	<i>21.36</i>	<i>1.345</i>	<i>-104.7</i>	<i>Olive green</i>	<i>Strong</i>
<i>1222</i>	<i>15</i>	<i>I</i>	<i>8.21</i>	<i>2.071</i>	—	<i>1.83</i>	<i>21.66</i>	<i>1.347</i>	<i>-109.5</i>	<i>Olive grey</i>	<i>Strong</i>
<i>1231</i>	<i>22</i>	<i>I</i>	<i>8.16</i>	<i>2.088</i>	—	<i>1.88</i>	<i>22.45</i>	<i>1.357</i>	<i>-109.8</i>	<i>Grey</i>	<i>Strong</i>
Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:				
<i>1</i>	<i>4-2-13</i>	<i>ice</i>	<i>8260B - VOCs + OxyS</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	<i>LL-15B-C40213 @ 12:44</i>				
<i>1</i>	<i>1244</i>	<i>ice</i>	<i>8015M - TPH-g</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>					

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(ft) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

GROUNDWATER SAMPLING LOG

Page 1 of 2

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013** *4-2-13*

WELL NO. **MW-15C** Hospital
 SAMPLED BY: **Frane Sosic**

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 2"	(inches)
DEPTH OF WELL <i>198.00</i>	(ft.)
DEPTH TO WATER <i>111.53</i>	(ft.)
HEIGHT OF WATER COLUMN <i>86.47</i>	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = <i>14.09461</i>	(gal)
PURGE VOLUME x 3 = <i>42.28383</i>	(gal)
PRODUCT THICKNESS	(ft.)

WELL CONDITION:

OK

WEATHER CONDITIONS:

Mostly sunny w/ light breeze

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp. Cond. <i>cm</i> S/cm	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
<i>1240</i>	<i>5</i>	<i>VAC TRUCK</i>	<i>8.14</i>	<i>1.959</i>	<i>—</i>	<i>3.21</i>	<i>23.33</i>	<i>1.269</i>	<i>107.6</i>	<i>light gray</i>	<i>Mild</i>
<i>1250</i>	<i>10</i>	<i>↓</i>	<i>8.05</i>	<i>1.950</i>	<i>—</i>	<i>2.79</i>	<i>24.06</i>	<i>1.260</i>	<i>105.3</i>	<i>light gray</i>	<i>Mild</i>
<i>1258</i>	<i>15</i>	<i>↓</i>	<i>8.02</i>	<i>1.952</i>	<i>—</i>	<i>2.13</i>	<i>23.10</i>	<i>1.265</i>	<i>102.1</i>	<i>cloudy</i>	<i>Mild</i>

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
<i>1</i>	<i>4-2-13</i>	<i>ice</i>	<i>8260B - VOCs + OxyS</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	<i>*Stinger must be raised ~15-18' to purge*</i> <i>LL-15C-040213 @ 14:36</i>
<i>1</i>	<i>14:36</i>	<i>ice</i>	<i>8015M - TPH-g</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

Page 2 of 2

PROJECT NO.: 1003-001-300

DATE: 4-2-13 (2013)

WELL NO. 15-C

SAMPLED BY: Frane Sosic

[illegible]

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013 4-3-13**

WELL NO. **MW-16A** Walker
 SAMPLED BY: **Frane Sasic**

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 2"	(inches)
DEPTH OF WELL 123.00	(ft.)
DEPTH TO WATER 112.37	(ft.)
HEIGHT OF WATER COLUMN 10.63	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = 1.73269	(gal)
PURGE VOLUME x 3 = 5.19807	(gal)
PRODUCT THICKNESS	(ft.)

WELL NOTES:
 WELL CONDITION: **GOOD**

WEATHER CONDITIONS:
Hazy AM & cool (~60°F)
Clears to sunny (~70°F)
 PURGING AND SAMPLING EQUIPMENT:
 YSI 556
 Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. $\mu\text{S/cm}$	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
1015	~3	VACUUM	7.92	2.698	--	3.86	20.96	1.754	-45.7	Olive green	Mild
1030	6	TRUCK	7.80	2.633	--	2.88	21.58	1.711	-59.6	Olive green	Mild

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	4-3-13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	Very slow purging well. Vacuum truck has a hose fitting leak, vacuum @ ~15 Hg"... optimal is ~20" Hg. LL-16A-040313 @ 10:47
1	1047	ice	8015M - TPH-g	VOAs	3	HCL	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

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GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013** *4-3-13*

WELL NO. **MW-16B** Walker
 SAMPLED BY: **Frane Sasic**

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 2"	(inches)
DEPTH OF WELL <i>160.00</i>	(ft.)
DEPTH TO WATER <i>114.92</i>	(ft.)
HEIGHT OF WATER COLUMN <i>42.08</i>	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = <i>6.85704</i>	(gal)
PURGE VOLUME x 3 = <i>20.5712</i>	(gal)
PRODUCT THICKNESS	(ft.)

WELL NOTES:
 WELL CONDITION: *GROD*

WEATHER CONDITIONS:
*Sunny w/ scattered clouds + light breeze
 (~78°F)*

PURGING AND SAMPLING EQUIPMENT:
 YSI 556
 Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. µS/cm	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
<i>1105</i>	<i>5</i>		<i>8.51</i>	<i>2.711</i>	<i>--</i>	<i>3.16</i>	<i>22.68</i>	<i>1.762</i>	<i>-140.2</i>	<i>Grey</i>	<i>Strong CH₄</i>
<i>1120</i>	<i>10</i>		<i>8.58</i>	<i>2.609</i>	<i>--</i>	<i>3.37</i>	<i>23.53</i>	<i>1.696</i>	<i>-169.5</i>	<i>Cloudy</i>	<i>Strong CH₄</i>
<i>1134</i>	<i>15</i>		<i>8.50</i>	<i>2.527</i>	<i>--</i>	<i>2.72</i>	<i>24.00</i>	<i>1.643</i>	<i>-192.4</i>	<i>Light grey</i>	<i>Strong CH₄</i>
<i>1144</i>	<i>20</i>		<i>8.45</i>	<i>2.496</i>	<i>--</i>	<i>2.10</i>	<i>24.35</i>	<i>1.622</i>	<i>-198.3</i>	<i>Cloudy</i>	<i>Strong CH₄</i>
Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:				
<i>1</i>	<i>4-3-13</i>	<i>ice</i>	<i>8260B - VOCs + Oxys</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	<i>Initial GW very slow to come up</i>				
<i>1</i>	<i>1203</i>	<i>ice</i>	<i>8015M - TPH-g</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	<i>LL-16B-040313 @ 1203</i>				

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

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GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013 4-3-13**

WELL NO. **MW-16C** Walker
 SAMPLED BY: **Frane Sasic**

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 2"	(inches)
DEPTH OF WELL 196.30	(ft.)
DEPTH TO WATER 117.70	(ft.)
HEIGHT OF WATER COLUMN 78.60	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = 12.8118	(gal)
PURGE VOLUME x 3 = 38.4354	(gal)
PRODUCT THICKNESS	(ft.)

WELL NOTES:
 WELL CONDITION:

GOOD

WEATHER CONDITIONS:
 Mostly sunny w/ light breeze
 (~74°F)

PURGING AND SAMPLING EQUIPMENT:
 YSI 556
 Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp. Cond. $\mu\text{S/cm}$	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
1303	5	VACUUM	8.23	1.491	--	2.55	24.98	0.968	-185.4	Dark grey	Strong Chlorine
1316	10	TRUCK	8.08	1.683	--	2.15	24.77	1.074	-147.9	Olive green	Strong Chlorine
1327	15		8.11	1.843	--	2.11	25.12	1.197	-166.3	Light green	Strong Chlorine

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	4-3-13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	Well would not pump at all really. After adjusting the stinger height and trying to get water to purge for quite some time, we ceased the purge to pull out and inspect the stinger in 16C. We found (2) cuts where there should only be (1) to allow air uptake that K segment was replaced and stinger was re-inserted into the well.
1	@	ice	8015M - TPH-g	VOAs	3	HCL	
1	1435	ice	8015 - Methane	VOAs	3	None	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

LL-16C_040313 @ 14:35

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PROJECT NO.: 1003-001-300

DATE: 4-3-2013 (202013)

WELL NO. 16-C

SAMPLED BY: Frane Sosic

[illegible]

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013** *4-9-13*

WELL NO. **W-17A** Lakeland
 SAMPLED BY: **Frane Sasic**

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 2"	(inches)
DEPTH OF WELL <i>108.30</i>	(ft.)
DEPTH TO WATER <i>97.25</i>	(ft.)
HEIGHT OF WATER COLUMN <i>11.05</i>	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = <i>1.8015</i>	(gal)
PURGE VOLUME x 3 = <i>5.4045</i>	(gal)
PRODUCT THICKNESS	(ft.)

WELL CONDITION:

OK

WEATHER CONDITIONS:

Clear/sunny/breezy (~70°F)

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. (S/cm)	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS	ORP	Color	Odor
<i>1008</i>	<i>5</i>	<i>VAC TRUCK</i>	<i>7.94</i>	<i>2.208</i>	<i>/</i>	<i>2.67</i>	<i>21.60</i>	<i>1.435</i>	<i>-27.8</i>	<i>Grey</i>	<i>Mild</i>

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES: <i>Very slow purge will de-aerated; allowed to recharge prior to sample collection: LL-17A-040913 @ 11:11</i>
<i>1</i>	<i>4-9-13</i>	<i>ice</i>	<i>8260B - VOCs + Oxys</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	
<i>1</i>	<i>11:11</i>	<i>ice</i>	<i>8015M - TPH-g</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

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GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013 4-9-13**

WELL NO. **W-17B** Lakeland
 SAMPLED BY: **Frane Sasic**

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 2"	(inches)
DEPTH OF WELL 169.60	(ft.)
DEPTH TO WATER 107.01	(ft.)
HEIGHT OF WATER COLUMN 62.59	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = 10.20217	(gal)
PURGE VOLUME x 3 = 30.60651	(gal)
PRODUCT THICKNESS	(ft.)

WELL CONDITION:

OK

WEATHER CONDITIONS:

Clear/sunny/breezy (~73°F)

PURGING AND SAMPLING EQUIPMENT:

YSI S56

Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. $\mu\text{S/cm}$	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
1143	5	WATERTRUCK	8.06	1.623	/	2.63	21.84	1.058	-97.3	Grey	Strong
1209	10	 	8.19	1.513	/	2.42	22.81	0.981	-105.9	Cloudy	Strong
1223	15	 	8.11	1.546	/	1.97	24.23	1.005	-146.3	Clear	Strong



Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES: Very slow purging deep well. LL-17B-040913 @ 1310
1	4-9-13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	
1	13:10	ice	8015M - TPH-g	VOAs	3	HCL	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2 h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

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PROJECT NO.: 1003-001-300

DATE: 4-9-2013 (202013)

WELL NO. 17-B

SAMPLED BY: Frane Sosic

[illegible]

GROUNDWATER SAMPLING LOG

PROJECT NAME: CENCO
PROJECT NO.: 1003-001-300
DATE: 2Q2013 4-9-13

WELL NO. W-17C Lakeland
SAMPLED BY: Frane Sasic

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 2"	(inches)
DEPTH OF WELL 200.00	(ft.)
DEPTH TO WATER 107.08	(ft.)
HEIGHT OF WATER COLUMN 92.92	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = 15.14596	(gal)
PURGE VOLUME x 3 = 45.43788	(gal)
PRODUCT THICKNESS	(ft.)

WELL CONDITION:

OK

WEATHER CONDITIONS:

Clear/sunny/breezy (+83°F)

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. (µS/cm)	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
1315	5	VAC TRUCK	8.33	1.335	—	2.67	23.24	0.868	-156.5	Cloudy	Strong
1327	10	↓	8.30	1.322	—	2.70	23.47	0.858	-157.4	Clear	Strong
1344	15	↓	8.27	1.317	—	2.56	23.71	0.855	-171.2	Clear	Strong

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES: Very slow purging deep well (TD = 200' logs) LL-17C_040913 @ 16:12
1	4-9-13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	
1	1612	ice	8015M - TPH-g	VOAs	3	HCL	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2 h (\text{ft}) \times 7.48 \text{ gal./ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

Page 2 of 2

PROJECT NO.: 1003-001-300

DATE: 4-9-2013 (202013)

WELL NO.

SAMPLED BY: Frane Sosic

[illegible]

GROUNDWATER SAMPLING LOG

PROJECT NAME: CENCO
 PROJECT NO.: 1003-001-300
 DATE: 2Q2013 4-5-13

WELL NO. MW-104A Lakeland
 SAMPLED BY: Frane Sosic

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL <u>100.00</u>	(ft.)
DEPTH TO WATER <u>92.84</u>	(ft.)
HEIGHT OF WATER COLUMN <u>7.16</u>	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = <u>4.7256</u>	(gal)
PURGE VOLUME <u>266</u> x 3 = <u>14.1768</u>	(gal)
PRODUCT THICKNESS	(ft.)

WELL CONDITION:
OK - lid will not secure

WEATHER CONDITIONS:
Scattered clouds w/ light wind (~67°F)

PURGING AND SAMPLING EQUIPMENT:
 YSI 556
 Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. (s/cm)	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS	ORP	Color	Odor
/	<u>5</u>	/	<u>* PURGED DRY *</u>	/	/	/	/	/	/	/	/
/	<u>10</u>	/		/	/	/	/	/	/	/	/
/	<u>15</u>	/		/	/	/	/	/	/	/	/

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
<u>1</u>	<u>4-5-13</u>	<u>ice</u>	<u>8260B - VOCs + Oxys</u>	<u>VOAs</u>	<u>3</u>	<u>HCL</u>	<u>Well 104-A purged dry @ ~6 gallons.</u> <u>No GW parameters could be measured.</u> <u>Well allowed to re-charge prior to collecting sample:</u> <u>LL-104A-040513 @ 1300</u>
<u>1</u>	<u>1300</u>	<u>ice</u>	<u>8015M - TPH-g</u>	<u>VOAs</u>	<u>3</u>	<u>HCL</u>	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2 h (\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013 4-4-13**

WELL NO. **MW-106A** Bloomfield
 SAMPLED BY: **Frane Sosis**

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL 110.00	(ft.)
DEPTH TO WATER 104.24	(ft.)
HEIGHT OF WATER COLUMN 5.76	(ft.)
CASING VOLUME* Hgt. x 0.163-Gal./Ft. = 3.8016	(gal)
PURGE VOLUME 0.66 x 3 = 11.4031	(gal)
PRODUCT THICKNESS	(ft.)

WELL NOTES:
 WELL CONDITION: OK - not great
 WEATHER CONDITIONS: Hazy morning (~65°F)
 PURGING AND SAMPLING EQUIPMENT:
 YSI 556
 Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. $\mu S/cm$	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
	5										
	10		* VOLUME *								

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	4-4-13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	8:27 106-A purged dry ~ 4 gal. (similar to previous sampling events.) Allowed to re-charge prior to sampling. LL 106A-040413 @ 1234
1	1234	ice	8015M - TPH-g	VOAs	3	HCL	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(ft) \times 7.48 \text{ gal./ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013 4-4-13**

WELL NO. **MW-107A** Bloomfield
 SAMPLED BY: **Frane Sosic**

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL 110.00	(ft.)
DEPTH TO WATER 104.03	(ft.)
HEIGHT OF WATER COLUMN 5.97	(ft.)
CASING VOLUME* Hgt. x 0.163 Gal./Ft. = 3.7402	(gal)
PURGE VOLUME 0.66 x 3 = 11.8206	(gal)
PRODUCT THICKNESS	(ft.)

WELL NOTES:
 WELL CONDITION: Very Good
 WEATHER CONDITIONS: Mostly sunny/humid/light breeze (~75°F)
 PURGING AND SAMPLING EQUIPMENT:
 YSI 556
 Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. (S/cm)	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS μ L	ORP mV	Color	Odor
1414	5	VAC TRUCK	8.34	1.913	—	3.01	25.72	1.244	-203.9	Cloudy	Strong CH ₄
1421	10	I	8.30	1.901	—	2.47	26.01	1.235	-210.4	Trans. green	Strong CH ₄
1430	15	I	8.28	1.874	—	2.55	26.40	1.218	-251.8	Trans. green	Strong CH ₄
1438	20	I	8.25	1.898	—	2.25	26.06	1.235	—	Trans. green	Strong CH ₄
Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:				
1	4-3-13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	VAC TRUCK HAD ISSUES. → pump would not work correctly Nieto swapped out truck #500 for #506. LL-107A-040413 @ 15:25				
1	1525	ice	8015M - TPH-g	VOAs	3	HCL					

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

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GROUNDWATER SAMPLING LOG

PROJECT NAME: CENCO
 PROJECT NO.: 1003-001-300
 DATE: 2Q2013 4-8-13

WELL NO. MW-503B Coaster
 SAMPLED BY: Frane Sasic

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL <u>108.67</u>	(ft.)
DEPTH TO WATER <u>101.60</u>	(ft.)
HEIGHT OF WATER COLUMN <u>7.07</u>	(ft.)
CASING VOLUME* Hgt. x 0.66 Gal./Ft. = <u>4.6662</u>	(gal)
PURGE VOLUME x 3 = <u>13.9986</u>	(gal)
PRODUCT THICKNESS	(ft.)

WELL NOTES:
 WELL CONDITION: Not good

WEATHER CONDITIONS:
Clear/sunny/gusty wind (~75°F)

PURGING AND SAMPLING EQUIPMENT:
 YSI 556
 Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. µS/cm	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
<u>1340</u>	<u>5</u>		<u>7.85</u>	<u>1.876</u>	-	<u>2.30</u>	<u>23.25</u>	<u>1.224</u>	<u>-2.5</u>	<u>Cloudy</u>	<u>Strong</u>
<u>1355</u>	<u>10</u>		<u>7.78</u>	<u>1.826</u>	-	<u>2.23</u>	<u>24.12</u>	<u>1.187</u>	<u>-17.2</u>	<u>Cloudy</u>	<u>Strong</u>
<u>1404</u>	<u>15</u>		<u>7.72</u>	<u>1.812</u>	-	<u>2.46</u>	<u>24.13</u>	<u>1.178</u>	<u>-31.5</u>	<u>Clear</u>	<u>Strong</u>

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
<u>1</u>	<u>4-8-13</u>	<u>ice</u>	<u>8260B - VOCs + Oxys</u>	<u>VOAs</u>	<u>3</u>	<u>HCL</u>	<u>LL-503B-040813-01 @ 1410</u> <u>LL-503B-040813-02 @ 1420</u>
<u>1</u>	<u>1410</u>	<u>ice</u>	<u>8015M - TPH-g</u>	<u>VOAs</u>	<u>3</u>	<u>HCL</u>	
<u>2</u>	<u>4-8-13</u>	<u>ice</u>	<u>8260B</u>	<u>VOAs</u>	<u>3</u>	<u>HCL</u>	
<u>2</u>	<u>1420</u>	<u>ice</u>	<u>8015M</u>	<u>VOAs</u>	<u>3</u>	<u>HCL</u>	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013 4-10-13**

WELL NO. **MW-701** Lakeland
 SAMPLED BY: **Frane Sosis**

Well Notes: _____

WELL CONDITION: _____

Very Good

WEATHER CONDITIONS: _____

Clear/sunny/light breeze (~85°F)

PURGING AND SAMPLING EQUIPMENT: _____

YSI 556

Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL <i>130.00</i>	(ft.)
DEPTH TO WATER <i>99.91</i>	(ft.)
HEIGHT OF WATER COLUMN <i>30.09</i>	(ft.)
CASING VOLUME* Hgt. x 0.66 Gal./Ft. = <i>19.8594</i>	(gal)
PURGE VOLUME x 3 = <i>59.5782</i>	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. $\mu S/cm$	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
<i>1312</i>	<i>5</i>	<i>VAC TRUCK</i>	<i>7.81</i>	<i>1.969</i>	<i>/</i>	<i>4.20</i>	<i>24.29</i>	<i>1.280</i>	<i>53.6</i>	<i>Light grey</i>	<i>None</i>
<i>1314</i>	<i>10</i>	<i>/</i>	<i>7.90</i>	<i>1.964</i>	<i>/</i>	<i>3.65</i>	<i>23.58</i>	<i>1.276</i>	<i>39.5</i>	<i>Light grey</i>	<i>Slight</i>
<i>1317</i>	<i>15</i>	<i>/</i>	<i>7.85</i>	<i>1.964</i>	<i>/</i>	<i>3.33</i>	<i>23.66</i>	<i>1.278</i>	<i>36.1</i>	<i>Light grey</i>	<i>Slight</i>

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
<i>1</i>	<i>4-10-13</i>	<i>ice</i>	<i>8260B - VOCs + Oxys</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	<i>LL-701-041013 @ 14:13</i>
<i>1</i>	<i>1413</i>	<i>ice</i>	<i>8015M - TPH-g</i>	<i>VOAs</i>	<i>3</i>	<i>HCL</i>	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2 h(ft) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

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CENCO

1003-001-300

4-10-13 (2Q2013)

70.

SAMPLED BY: Frane Sosic

[illegible]

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013 4-10-13**

WELL NO. **MW-702** Lakeland
 SAMPLED BY: **Frane Sosis**

Well Notes: **Strong H₂S / CH₄ / VOC vapors**

WELL CONDITION:

WEATHER CONDITIONS:

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL 130.00	(ft.)
DEPTH TO WATER 99.52	(ft.)
HEIGHT OF WATER COLUMN 30.48	(ft.)
CASING VOLUME* Hgt. x .66 Gal./Ft. = 20.1168	(gal)
PURGE VOLUME x 3 = 60.3504	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. (µS/cm)	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
1452	5	VAC TRUCK	7.84	2.198	—	3.15	26.39	1.427	19.1	Clear	Strong
1453	10	↓	7.77	2.184	—	3.01	25.87	1.420	-18.6	Clear	Strong
1455	15	↓	7.68	2.184	—	2.97	25.99	1.419	-29.7	Clear	Strong

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES: Vent well for 4+ hours prior to sampling
1	4-10-13	ice	8260B - VOCs + OxyS	VOAs	3	HCL	LL-702_041013 @ 15:42
1	@	ice	8015M - TPH-g	VOAs	3	HCL	
1	1542	ice	8015 - Methane	VOAs	3	NONE	
1	-11-	ice	Total Alkalinity	250 ml poly	1	NONE	
1	-11-	ice	300 IC: nitrate, sulfate	250 ml poly	1	NONE	
1	-11-	ice	Ferrous Iron	250 ml poly	1	HNO ₃	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

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PROJECT NO.: 1003-001-300

DATE: 4-10-13 (202013)

702

SAMPLED BY: Frane Sosic

[illegible]

GROUNDWATER SAMPLING LOG

PROJECT NAME: CENCO
PROJECT NO.: 1003-001-300
DATE: 2Q2013 4-10-13

WELL NO. MW-703 Lakeland
SAMPLED BY: Frane Sosic

Well Notes: Strong H2S / LEL / VOC vapors

WELL CONDITION:

GOOD

WEATHER CONDITIONS:

Clear / sunny / light breeze (~85°F)

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL 130.00	(ft.)
DEPTH TO WATER 101.17	(ft.)
HEIGHT OF WATER COLUMN 28.83	(ft.)
CASING VOLUME* Hgt. x 0.66 Gal./Ft. = 19.0278	(gal)
PURGE VOLUME x 3 = 57.0834	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. (s/cm)	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS μ C	ORP mV	Color	Odor
1210	5	UP TRACK	7.98	1.944	—	4.35	23.54	1.263	-3.5	Light grey	Strong
1213	10		7.92	1.949	—	3.63	23.32	1.267	-36.1	Light grey	Strong
1215	15		7.91	1.951	—	3.15	23.31	1.268	-63.6	Light grey	Strong

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	4-10-13	ice	8260B - VOCs + OxyS	VOAs	3	HCL	LL_703_041013 @ 12:45
1	@	ice	8015M - TPH-g	VOAs	3	HCL	
1	1245	ice	8015 - Methane	VOAs	3	NONE	
1	-11-	ice	Total Alkalinity	250 ml poly	1	NONE	
1	-11-	ice	300 IC: nitrate, sulfate	250 ml poly	1	NONE	
1	-11-	ice	Ferrous Iron	250 ml poly	1	HNO ₃	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal./ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

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PROJECT NO.: 1003-001-300

DATE: 4-10-13 (2020/3)

WELL NO.

SAMPLED BY: Frane Sosic

[illegible]

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **04/15/2013 2Q2013**

WELL NO. **MW-704** Lakeland
 SAMPLED BY: **Frank Sasic AW**

Well Notes: **Strong H2S / LEL / VOC vapors**

WELL CONDITION:

GOOD

WEATHER CONDITIONS:

OVERCAST, COOL

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL 130.00	(ft.)
DEPTH TO WATER 102.90	(ft.)
HEIGHT OF WATER COLUMN 27.1	(ft.)
CASING VOLUME* Hgt. x 0.66 Gal./Ft. = 17.9	(gal)
PURGE VOLUME x 3 = 53.7	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. (Ms/cm)	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
1014	5	VALTRUCK	7.94	2.135	/	2.05	21.13	1.389	-121.7	CLEAR	SLIGHT
1016	10	↓	7.90	2.134	/	2.07	22.41	1.388	-132.7	CLEAR	SLIGHT
1018	15	↓	7.78	2.152	/	2.09	22.59	1.391	-143.6	CLEAR	SLIGHT

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	04/15/13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	LL - 704 - 041513 @ 1130 VACUUM/RECHARGE VERY SLOW AFTER PURGE 30 GAL. SAMPLED AT PURGE 45 GAL.
1	↻	ice	8015M - TPH-g	VOAs	3	HCL	
1	↻	ice	8015 - Methane	VOAs	3	NONE	
1	1130	ice	Total Alkalinity	250 ml poly	1	NONE	
1		ice	300 IC: nitrate, sulfate	250 ml poly	1	NONE	
1		ice	Ferrous Iron	250 ml poly	1	HNO ₃	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2 h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

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PROJECT NO.: 1003-001-300

DATE: 04/15/13 (2Q2013)

SAMPLED BY: ~~Frane Sasic~~ AW

[illegible]

GROUNDWATER SAMPLING LOG

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PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013 4-10-13**

WELL NO. **MW-705** Lakeland
 SAMPLED BY: **Frane Sosic + AW**

Well Notes: **Strong H2S / LEL / VOC vapors**

WELL CONDITION: **GOOD**

WEATHER CONDITIONS:

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL 130.00	(ft.)
DEPTH TO WATER 103.90	(ft.)
HEIGHT OF WATER COLUMN 26.10	(ft.)
CASING VOLUME* Hgt. x 0.66 Gal./Ft. = 17.23	(gal)
PURGE VOLUME x 3 = 51.68	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. $\mu S/cm$	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP μV	Color	Odor
1602	5	WATER TRUCK	7.65	2.060	/	2.67	24.27	1.337	67.2	Cloudy	Strong
1604	10	↓	7.77	2.055	/	3.09	24.62	1.336	58.6	-11-	-11-
1606	20	↓	7.77	2.056	/	2.82	24.57	1.336	47.0	Cloudy	Strong

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES: Vent well for 4+ hours prior to sampling
1	4-10-13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	LL-705-041013 @ 16:36
1	@	ice	8015M - TPH-g	VOAs	3	HCL	
1	16:36	ice	8015 - Methane	VOAs	3	NONE	
1	-11-	ice	Total Alkalinity	250 ml poly	1	NONE	
1	-11-	ice	300 IC: nitrate, sulfate	250 ml poly	1	NONE	
1	-11-	ice	Ferrous Iron	250 ml poly	1	HNO₃	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(ft) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

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Page 2 of 2

PROJECT NO.: 1003-001-300

DATE: 4-10-13 (2Q2013)

WELL NO.

SAMPLED BY: Frane Sosic + AW

[illegible]

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **04/15/2013 2Q2013**

WELL NO. **MW-706** Lakeland
 SAMPLED BY: **Frank Sosie AW**

Well Notes: _____
 WELL CONDITION: **GOOD**

WEATHER CONDITIONS: **OVERCAST, LIGHT RAIN**

PURGING AND SAMPLING EQUIPMENT:
 YSI 556
 Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL 130.00	(ft.)
DEPTH TO WATER 100.48	(ft.)
HEIGHT OF WATER COLUMN 29.52	(ft.)
CASING VOLUME* Hgt. x 0.46 Gal./Ft. = 19.5	(gal)
PURGE VOLUME x 3 = 58.4	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp. Cond. $\mu\text{S/cm}$	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
0846	5	VAC TRUCK	8.04	2.024	/	2.72	21.28	1.315	-68.4	CLEAR	STRONG
0848	10	↓	8.05	2.038	/	2.71	22.82	1.322	-90.7	CLEAR	STRONG
0850	15	↓	8.13	2.035	/	2.38	22.73	1.322	-107.3	CLEAR	STRONG

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1/2	04/15/13	ice	8260B - VOCs + Oxys	VOAs	3/3	HCL	LL-706-041513-01 @ 0920 LL-706-041513-02 @ 0930
1/2	C	ice	8015M - TPH-g	VOAs	3/3	HCL	
1		ice	8015 - Methane	VOAs	3	NONE	
1	0920(1)	ice	Total Alkalinity	250 ml poly	1	NONE	
1	0930(2)	ice	300 IC: nitrate, sulfate	250 ml poly	1	NONE	
1		ice	Ferrous Iron	250 ml poly	1	HNO ₃	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2 h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

Page 2 of 2

PROJECT NO.: 1003-001-300

DATE: 04/15/13 (2Q 2013)

SAMPLED BY: ~~Frane Sosic~~ AW

[illegible]

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **2Q2013 4-8-13**

WELL NO. **MW-707** Coaster
 SAMPLED BY: **Frane Sosic**

Well Notes: _____
 WELL CONDITION: **GOOD**

WEATHER CONDITIONS: **Clear/sunny/gusty winds (~75°F)**

PURGING AND SAMPLING EQUIPMENT:
 YSI 556
 Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL 130.00	(ft.)
DEPTH TO WATER 98.46	(ft.)
HEIGHT OF WATER COLUMN 31.54	(ft.)
CASING VOLUME* Hgt. x 0.66 Gal./Ft. = 20.8164	(gal)
PURGE VOLUME x 3 = 62.4492	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. (s/cm)	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS	ORP mV	Color	Odor
1440	5		7.71	1.764	--	5.03	22.92	1.148	-3.5	Light grey	Strong
1442	10		7.69	1.771	--	3.10	22.84	1.154	-15.8	Light grey	Strong
1445	15		7.66	1.787	--	2.59	22.60	1.162	-23.6	Light grey	Strong

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	4-8-13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	LL-707-040813 @ 1543
1	15:43	ice	8015M - TPH-g	VOAs	3	HCL	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

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PROJECT NO.: 1003-001-300

DATE: 4-8-2013 (2020/3)

WELL NO. 701

SAMPLED BY: Frane Sosic

[illegible]

GROUNDWATER SAMPLING LOG

PAGE 1 OF 2

PROJECT NAME: **CENCO**
PROJECT NO.: **1003-001-300**
DATE: **04/11/2013 2Q2013**

WELL NO. **MW-708** Hospital
SAMPLED BY: **Frane Sosic AW**

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL 130.00	(ft.)
DEPTH TO WATER FPPH 97.55	(ft.)
HEIGHT OF WATER COLUMN 32.45	(ft.)
CASING VOLUME* Hgt. x 0.66 Gal./Ft. = 21.4	(gal)
PURGE VOLUME x 3 = 64.3	(gal)
PRODUCT THICKNESS 97.82 (DTW) - 97.55 (FP) = 0.27 ON	(ft.)

Well Notes: **May contain FPPH**

WELL CONDITION:
GOOD

WEATHER CONDITIONS:
WARM, SUNNY

PURGING AND SAMPLING EQUIPMENT:
YSI 556
Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. $\mu\text{S}/\text{cm}$	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
-	5	VACTRUK	* PARAMETERS NOT TAKEN							GRAY	MODERATE
-	10		DUE TO FPPH (CO.5 GAL)							LT. GRAY	SLIGHT
1508	15	↓	7.75	2.063		2.42	23.27	1.340	-23.2	LT. GRAY	SLIGHT

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1/2	04/11/13	ice	8260B - VOCs + Oxys	VOAs	3/3	HCL	* 30 FT OF STINGER REMOVED TO SKIM FIRST 10 GAL WATER, THEN LOWERED AND CONTINUED AS NORMAL LL-708-041113-01 @ 1615 LL-708-041113-02 @ 1620
1/2	e	ice	8015M - TPH-g	VOAs	3/3	HCL	
1		ice	8015 - Methane	VOAs	3	NONE	
1	1615 (1)	ice	Total Alkalinity	250 ml poly	1	NONE	
1	1620 (2)	ice	300 IC: nitrate, sulfate	250 ml poly	1	NONE	
1		ice	Ferrous Iron	250 ml poly	1	HNO ₃	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2 h(\text{ft}) \times 7.48 \text{ gal}/\text{ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

Page 2 of 2

PROJECT NO.: 1003-001-300

DATE: 04/11/13 (2Q 2013)

WELL NO. 708

SAMPLED BY: ~~Frank Sosic~~ AW

[illegible]

GROUNDWATER SAMPLING LOG

PAGE 1 of 2

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **04/11/2013 2Q2013**

WELL NO. **MW-709** Hospital
 SAMPLED BY: **Frank Sasic AW**

Well Notes: _____

WELL CONDITION: _____

GOOD

WEATHER CONDITIONS: _____

WARM, SUNNY

PURGING AND SAMPLING EQUIPMENT: _____

YSI 556

Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL <u>130.00</u>	(ft.)
DEPTH TO WATER <u>109.65</u>	(ft.)
HEIGHT OF WATER COLUMN <u>20.35</u>	(ft.)
CASING VOLUME* Hgt. x 0.66 Gal./Ft. = <u>13.4</u>	(gal)
PURGE VOLUME x 3 = <u>40.3</u>	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. uS/cm	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
1309	5	VALTRUCK	7.80	2.207	/	2.71	22.97	1.435	84.9	CLOUDY	SLIGHT
1312	10	↓	7.76	2.206	/	2.50	23.45	1.434	59.3	CLOUDY	SLIGHT
1316	15	↓	7.80	2.213	/	2.61	23.40	1.442	48.3	CLOUDY	SLIGHT

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	04/11/13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	LL-709-041113 @ 1350
1	1350	ice	8015M - TPH-g	VOAs	3	HCL	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

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PROJECT NO.: 1003-001-300

DATE: 04/11/13 (2Q 2013)

WELL NO. 709

SAMPLED BY: ~~Frane Sosic~~ AW

[illegible]

GROUNDWATER SAMPLING LOG

PAGE 1 of 2

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **04/12/2013 2Q2013**

WELL NO. **MW-710** Hospital
 SAMPLED BY: **Frane Sasic AW**

Well Notes: _____

WELL CONDITION: _____

GOOD

WEATHER CONDITIONS: _____

OVERCAST, COOL

PURGING AND SAMPLING EQUIPMENT: _____

YSI 556

Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL 130.00	(ft.)
DEPTH TO WATER 96.25	(ft.)
HEIGHT OF WATER COLUMN 33.75	(ft.)
CASING VOLUME* Hgt. x 0.66 Gal./Ft. = 22.3	(gal)
PURGE VOLUME x 3 = 66.8	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. μS/cm	Turbidity NTUs	DO mg/L	Temperature (F/°C)	TDS g/L	ORP mV	Color	Odor
0901	5	VAR TRUCK	7.95	1.843	/	2.75	21.35	1.198	66.7	cloudy	—
0903	10	↓	7.89	1.844	/	2.71	21.39	1.198	62.3	CLOUDY	—
0905	15	↓	7.83	1.847	/	2.59	21.78	1.200	59.6	CLOUDY	—

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	04/12/13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	LL - 710 - 041213 @ 0925
1	e	ice	8015M - TPH-g	VOAs	3	HCL	
	0925						

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2 h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

Page 2 of 2

PROJECT NO.: 1003-001-300

DATE: 04/12/13 (20 2013)

SAMPLED BY: ~~Frane Sosic~~ AW

[illegible]

GROUNDWATER SAMPLING LOG

PAGE 1 of 2

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **04/12/2013 2Q2013**

WELL NO. **MW-711** Hospital
 SAMPLED BY: ~~Franc Sosie~~ **AW**

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL 130.00	(ft.)
DEPTH TO WATER 103.00	(ft.)
HEIGHT OF WATER COLUMN 27.00	(ft.)
CASING VOLUME* Hgt. x 0.66 Gal./Ft. = 17.8	(gal)
PURGE VOLUME x 3 = 53.5	(gal)
PRODUCT THICKNESS	(ft.)

Well Notes:

WELL CONDITION:

GOOD

WEATHER CONDITIONS:

OVERCAST, COOL

PURGING AND SAMPLING EQUIPMENT:

YSI 556

Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. μ S/cm	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
0943	5	VAL TRUCK	7.70	1.764	/	2.62	21.43	1.145	-45.4	DK GRAY	STRONG
0946	10	↓	7.57	1.765	/	2.42	22.13	1.146	-82.5	DK GRAY	STRONG
0949	15	↓	7.64	1.782	/	1.96	22.20	1.157	-83.3	GRAY	STRONG

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	04/12/13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	LL - 711 - 041213 @ 1020
1	@	ice	8015M - TPH-g	VOAs	3	HCL	
1	1020	ice	8015 - Methane	VOAs	3	NONE	
1		ice	Total Alkalinity	250 ml poly	1	NONE	
1		ice	300 IC: nitrate, sulfate	250 ml poly	1	NONE	
1		ice	Ferrous Iron	250 ml poly	1	HNO ₃	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2 h (ft) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

[illegible]

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **04/12/2013 2Q2013**

WELL NO. **MW-712** Hospital
 SAMPLED BY: **Frank Sesie Aw**

Well Notes: _____

WELL CONDITION: _____

GOOD

WEATHER CONDITIONS: _____

SUNNY, WARM

PURGING AND SAMPLING EQUIPMENT: _____

YSI 556

Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL 130.00	(ft.)
DEPTH TO WATER 99.95	(ft.)
HEIGHT OF WATER COLUMN 30.05	(ft.)
CASING VOLUME* Hgt. x 0.66 Gal./Ft. = 19.8	(gal)
PURGE VOLUME x 3 = 59.5	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. μ S/cm	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
1347	5	VALTRUCK	7.67	1.806	/	2.79	24.58	1.174	-69.4	CLOUDY	SLIGHT
1349	10	↓	7.59	1.807	/	2.51	24.17	1.175	-97.7	CLOUDY	SLIGHT
1352	15	↓	7.62	1.806	/	2.39	24.01	1.173	-110.4	CLOUDY	SLIGHT

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	04/12/13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	LL-712-041213 @ 1420
1	1420	ice	8015M - TPH-g	VOAs	3	HCL	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

Page 12 of 12

WELL NO. 712

SAMPLED BY: ~~Frane Sosic~~ AW

DATE: 04/12/13 (2013)

[illegible]

GROUNDWATER SAMPLING LOG

PAGE 1 of 2

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **04/11/2013** **2Q2013**

WELL NO. **MW-713** Hospital
 SAMPLED BY: **Erane Sasic AW**

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL 130.00	(ft.)
DEPTH TO WATER 105.59	(ft.)
HEIGHT OF WATER COLUMN 24.41	(ft.)
CASING VOLUME* Hgt. x 0.66 Gal./Ft. = 16.1	(gal)
PURGE VOLUME x 3 = 48.3	(gal)
PRODUCT THICKNESS	(ft.)

Well Notes: _____
 WELL CONDITION: _____

GOOD

WEATHER CONDITIONS: _____
 WARM, SUNNY

PURGING AND SAMPLING EQUIPMENT: _____
 YSI 556
 Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. μ S/cm	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
1028	5	VAC TRACK	7.89	2.065	/	2.95	23.29	1.343	9.1	CLOUDY	SLIGHT
1030	10	↓	7.82	2.079	/	2.89	23.73	1.347	-0.5	CLOUDY	SLIGHT
1033	15	↓	7.67	2.189	/	2.81	23.92	1.421	-88.6	CLOUDY	SLIGHT

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1/2	04/11/13	ice	8260B - VOCs + Oxys	VOAs	3/3	HCL	LL-713-041113-01 C 1055 LL-713-041113-02 C 1105
1/2	@	ice	8015M - TPH-g	VOAs	3/3	HCL	
1	@	ice	8015 - Methane	VOAs	3	NONE	
1	1055 (1)	ice	Total Alkalinity	250 ml poly	1	NONE	
1	1105 (2)	ice	300 IC: nitrate, sulfate	250 ml poly	1	NONE	
1		ice	Ferrous Iron	250 ml poly	1	HNO ₃	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2 h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

Page 2 of 2

PROJECT NO.: 1003-001-300

DATE: 04/11/13 (2Q 2013)

WELL NO. 713

SAMPLED BY: ~~Frank Sosie~~ AW

[illegible]

GROUNDWATER SAMPLING LOG

PAGE 1 of 2

PROJECT NAME: CENCO
 PROJECT NO.: 1003-001-300
 DATE: 04/11/2013 2Q2013

WELL NO. MW-714 Hospital
 SAMPLED BY: Frane Sosic AW

Well Notes: _____

WELL CONDITION: _____

GOOD

WEATHER CONDITIONS: _____

WARM, SUNNY

PURGING AND SAMPLING EQUIPMENT: _____

YSI 556

Interface probe (200')

WELL INFORMATION	
TOP OF CASING ELEV.	(ft.)
WELL DIAMETER 4"	(inches)
DEPTH OF WELL 142.00	(ft.)
DEPTH TO WATER 106.18	(ft.)
HEIGHT OF WATER COLUMN 35.82	(ft.)
CASING VOLUME* Hgt. x 0.66 Gal./Ft. = 23.6	(gal)
PURGE VOLUME x 3 = 70.9	(gal)
PRODUCT THICKNESS	(ft.)

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. μs/cm	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
0927	5	VAC TRAIL	7.95	2.509	/	4.30	22.22	1.631	3.7	CLOUDY	SLIGHT
0929	10	↓	7.85	2.521	/	3.34	22.55	1.634	-29.1	CLEAR	SLIGHT
0931	15	↓	7.81	2.511	/	2.91	22.92	1.633	-39.5	CLEAR	SLIGHT

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	04/11/13	ice	8260B - VOCs + OxyS	VOAs	3	HCL	LL-714-041113 @ 1005
1	@ 1005	ice	8015M - TPH-g	VOAs	3	HCL	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

Page 2 of 2

PROJECT NO.: 1003-001-300

DATE: 04/11/13 (2Q2013)

WELL NO. 714

SAMPLED BY: ~~Frane Sosic~~ AW

[illegible]

GROUNDWATER SAMPLING LOG

PROJECT NAME: **CENCO**
 PROJECT NO.: **1003-001-300**
 DATE: **04/12/2013 2Q2013**

WELL NO. **MW-715** Hospital
 SAMPLED BY: **Frane Sasic AW**

WELL INFORMATION		
TOP OF CASING ELEV.		(ft.)
WELL DIAMETER	4"	(inches)
DEPTH OF WELL	134.00	(ft.)
DEPTH TO WATER	97.98	(ft.)
HEIGHT OF WATER COLUMN	36.02	(ft.)
CASING VOLUME*	Hgt. x 0.66 Gal./Ft. = 23.8	(gal)
PURGE VOLUME	x 3 = 71.3	(gal)
PRODUCT THICKNESS		(ft.)

Well Notes: _____
 WELL CONDITION: _____

GOOD

WEATHER CONDITIONS: _____
SUNNY, WARM

PURGING AND SAMPLING EQUIPMENT: _____
 YSI 556
 Interface probe (200')

PURGE DATA											
Time:	Purge Volume (Gal.)	Flow Rate (Gal./Min.)	pH	Sp.Cond. μs/cm	Turbidity NTUs	DO mg/L	Temperature (F/C)	TDS g/L	ORP mV	Color	Odor
1124	5	✓	7.87	1.570	/	3.36	22.63	1.020	-99.2	CLOUDY	SLIGHT
1127	10	✓	7.80	1.567	/	2.91	22.94	1.018	-106.3	CLOUDY	SLIGHT
1131	15	✓	7.81	1.567	/	2.70	23.25	1.017	-121.5	CLOUDY	SLIGHT

Sample No.	Sample Time	Packing	Analyses	Container	Quantity	Preservative	NOTES:
1	04/12/13	ice	8260B - VOCs + Oxys	VOAs	3	HCL	LL - 715 - 041213 @ 1215
1	0	ice	8015M - TPH-g	VOAs	3	HCL	
1	1215	ice	8015 - Methane	VOAs	3	NONE	
1		ice	Total Alkalinity	250 ml poly	1	NONE	
1		ice	300 IC: nitrate, sulfate	250 ml poly	1	NONE	
1		ice	Ferrous Iron	250 ml poly	1	HNO ₃	

ADDITIONAL INFORMATION:

TOC = Top of well casing

*Casing Volume = $r^2 h(\text{ft}) \times 7.48 \text{ gal/ft.}^3$

4" well = 0.66 Gal./Foot

2" well = 0.163 Gal./Foot

Page 2 of 2

PROJECT NO.: 1003-001-300

DATE: 04/12/13 (202013)

715

SAMPLED BY: ~~Frane Sosic~~ AW

[illegible]

Appendix B



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

09 April 2013

Jeremy Squire
Murex
15375 Barranca Parkway, Suite K-101
Irvine, CA 92861
RE: Cenco

Enclosed are the results of analyses for samples received by the laboratory on 04/01/13 16:35. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Katherine Shields
Jr. Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
LL_14A_040113	T130756-01	Water	04/01/13 10:45	04/01/13 16:35
LL_14B_040113	T130756-02	Water	04/01/13 12:48	04/01/13 16:35
LL_14C_040113	T130756-03	Water	04/01/13 15:56	04/01/13 16:35
LL_TB_040113	T130756-04	Water	04/01/13 00:00	04/01/13 16:35
LL_W10_040113	T130756-05	Water	04/01/13 16:10	04/01/13 16:35

SunStar Laboratories, Inc.

Katherine Shields

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Katherine Shields, Jr. Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

LL_14A_040113
T130756-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	96	50	ug/l	1	3040311	04/03/13	04/04/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		95.5 %	65-135		"	"	"	"	

Metals by SM 3500 Series Methods

Ferrous Iron	ND	0.100	mg/l	1	3040247	04/02/13	04/08/13	EPA6010/S M3500	
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Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040222	04/02/13	04/03/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

LL_14A_040113
T130756-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1-Dichloroethene	ND	1.0	ug/l	1	3040222	04/02/13	04/03/13	EPA 8260B
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	91	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	2.4	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	12	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	5.8	0.50	"	"	"	"	"	"
Toluene	1.8	0.50	"	"	"	"	"	"
Ethylbenzene	1.4	0.50	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

LL_14A_040113
T130756-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

m,p-Xylene	6.6	1.0	ug/l	1	3040222	04/02/13	04/03/13	EPA 8260B	
o-Xylene	4.5	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	ND	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"	
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene		115 %	83.5-119		"	"	"	"	
Surrogate: Dibromofluoromethane		106 %	81-136		"	"	"	"	
Surrogate: Toluene-d8		99.6 %	88.8-117		"	"	"	"	

Conventional Chemistry Parameters by APHA/EPA/ASTM Methods

Total Alkalinity	460	20	mg/l	1	3040231	04/02/13	04/02/13	EPA 310.1	
Anions by EPA Method 300.0									
Sulfate as SO4	183	5.00	mg/l	10	3040312	04/03/13	04/03/13	EPA 300.0	
Nitrate as NO3	2.99	0.500	"	1	3040228	04/02/13	04/02/13	"	
RSK-175									
Methane	ND	1.00	ug/l	1	3040216	04/02/13	04/02/13	RSK-175	

SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

LL_14B_040113
T130756-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	ND	50	ug/l	1	3040311	04/03/13	04/04/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		94.0 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040222	04/02/13	04/03/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields, Jr. Project Manager



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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

LL_14B_040113
T130756-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040222	04/02/13	04/03/13	EPA 8260B	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	ND	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	21	1.0	"	"	"	"	"	"	
n-Propylbenzene	ND	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	2.6	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	ND	0.50	"	"	"	"	"	"	
Toluene	0.61	0.50	"	"	"	"	"	"	
Ethylbenzene	ND	0.50	"	"	"	"	"	"	
m,p-Xylene	2.0	1.0	"	"	"	"	"	"	
o-Xylene	1.2	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

LL_14B_040113
T130756-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3040222	04/02/13	04/03/13	EPA 8260B
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		111 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		105 %	81-136		"	"	"	"
Surrogate: Toluene-d8		99.6 %	88.8-117		"	"	"	"

SunStar Laboratories, Inc.

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/09/13 09:44
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LL_14C_040113
T130756-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	ND	50	ug/l	1	3040311	04/03/13	04/04/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		92.5 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040222	04/02/13	04/03/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

LL_14C_040113
T130756-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040222	04/02/13	04/03/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	27	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	2.4	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	0.88	0.50	"	"	"	"	"	"
Ethylbenzene	0.58	0.50	"	"	"	"	"	"
m,p-Xylene	2.7	1.0	"	"	"	"	"	"
o-Xylene	1.7	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

LL_14C_040113
T130756-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3040222	04/02/13	04/03/13	EPA 8260B	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"	
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene		113 %	83.5-119		"	"	"	"	
Surrogate: Dibromofluoromethane		110 %	81-136		"	"	"	"	
Surrogate: Toluene-d8		98.9 %	88.8-117		"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

LL_TB_040113
T130756-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040222	04/02/13	04/03/13	EPA 8260B
Bromochloromethane	ND	1.0	"	"	"	"	"	"
Bromodichloromethane	ND	1.0	"	"	"	"	"	"
Bromoform	ND	1.0	"	"	"	"	"	"
Bromomethane	ND	1.0	"	"	"	"	"	"
n-Butylbenzene	ND	1.0	"	"	"	"	"	"
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"
Chlorobenzene	ND	1.0	"	"	"	"	"	"
Chloroethane	ND	1.0	"	"	"	"	"	"
Chloroform	ND	1.0	"	"	"	"	"	"
Chloromethane	ND	1.0	"	"	"	"	"	"
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"
Dibromochloromethane	ND	1.0	"	"	"	"	"	"
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"
Dibromomethane	ND	1.0	"	"	"	"	"	"
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

LL_TB_040113
T130756-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

cis-1,3-Dichloropropene	ND	0.50	ug/l	1	3040222	04/02/13	04/03/13	EPA 8260B	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	ND	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	ND	1.0	"	"	"	"	"	"	
n-Propylbenzene	ND	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	ND	0.50	"	"	"	"	"	"	
Toluene	ND	0.50	"	"	"	"	"	"	
Ethylbenzene	ND	0.50	"	"	"	"	"	"	
m,p-Xylene	ND	1.0	"	"	"	"	"	"	
o-Xylene	ND	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	ND	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

LL_TB_040113
T130756-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	ug/l	1	3040222	04/02/13	04/03/13	EPA 8260B	
Surrogate: 4-Bromofluorobenzene	108 %	83.5-119			"	"	"	"	
Surrogate: Dibromofluoromethane	107 %	81-136			"	"	"	"	
Surrogate: Toluene-d8	98.2 %	88.8-117			"	"	"	"	

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Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

LL_W10_040113
T130756-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	490	50	ug/l	1	3040311	04/03/13	04/04/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene	96.1 %	65-135	"	"	"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040222	04/02/13	04/03/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

LL_W10_040113
T130756-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040222	04/02/13	04/03/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	13	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	6.3	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

LL_W10_040113
T130756-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	150	10	ug/l	1	3040222	04/02/13	04/03/13	EPA 8260B	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"	
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene		108 %	83.5-119		"	"	"	"	
Surrogate: Dibromofluoromethane		109 %	81-136		"	"	"	"	
Surrogate: Toluene-d8		98.1 %	88.8-117		"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

Purgeable Petroleum Hydrocarbons by EPA 8015C - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040311 - EPA 5030 GC

Blank (3040311-BLK1)

Prepared: 04/03/13 Analyzed: 04/04/13

C6-C12 (GRO)	ND	50	ug/l							
Surrogate 4-Bromofluorobenzene	95.2		"	100		95.2	65-135			

LCS (3040311-BS1)

Prepared: 04/03/13 Analyzed: 04/04/13

C6-C12 (GRO)	4930	50	ug/l	5520		89.2	75-125			
Surrogate 4-Bromofluorobenzene	83.7		"	100		83.7	65-135			

LCS Dup (3040311-BSD1)

Prepared: 04/03/13 Analyzed: 04/04/13

C6-C12 (GRO)	5430	50	ug/l	5520		98.4	75-125	9.74	20	
Surrogate 4-Bromofluorobenzene	70.7		"	100		70.7	65-135			

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

Metals by SM 3500 Series Methods - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040247 - EPA 3010A

Blank (3040247-BLK1)

Prepared: 04/02/13 Analyzed: 04/08/13

Ferrous Iron	ND	0.100	mg/l							
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Duplicate (3040247-DUP1)

Source: T130772-01

Prepared: 04/02/13 Analyzed: 04/08/13

Ferrous Iron	0.326	0.100	mg/l		0.383			16.0	200	
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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040222 - EPA 5030 GCMS

Blank (3040222-BLK1)

Prepared & Analyzed: 04/02/13

Bromobenzene	ND	1.0	ug/l
Bromochloromethane	ND	1.0	"
Bromodichloromethane	ND	1.0	"
Bromoform	ND	1.0	"
Bromomethane	ND	1.0	"
n-Butylbenzene	ND	1.0	"
sec-Butylbenzene	ND	1.0	"
tert-Butylbenzene	ND	1.0	"
Carbon tetrachloride	ND	0.50	"
Chlorobenzene	ND	1.0	"
Chloroethane	ND	1.0	"
Chloroform	ND	1.0	"
Chloromethane	ND	1.0	"
2-Chlorotoluene	ND	1.0	"
4-Chlorotoluene	ND	1.0	"
Dibromochloromethane	ND	1.0	"
1,2-Dibromo-3-chloropropane	ND	1.0	"
1,2-Dibromoethane (EDB)	ND	1.0	"
Dibromomethane	ND	1.0	"
1,2-Dichlorobenzene	ND	1.0	"
1,3-Dichlorobenzene	ND	1.0	"
1,4-Dichlorobenzene	ND	1.0	"
Dichlorodifluoromethane	ND	0.50	"
1,1-Dichloroethane	ND	1.0	"
1,2-Dichloroethane	ND	0.50	"
1,1-Dichloroethene	ND	1.0	"
cis-1,2-Dichloroethene	ND	1.0	"
trans-1,2-Dichloroethene	ND	1.0	"
1,2-Dichloropropane	ND	1.0	"
1,3-Dichloropropane	ND	1.0	"
2,2-Dichloropropane	ND	1.0	"
1,1-Dichloropropene	ND	1.0	"
cis-1,3-Dichloropropene	ND	0.50	"
trans-1,3-Dichloropropene	ND	0.50	"
Hexachlorobutadiene	ND	1.0	"
Isopropylbenzene	ND	1.0	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

Volatil Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040222 - EPA 5030 GCMS

Blank (3040222-BLK1)

Prepared & Analyzed: 04/02/13

p-Isopropyltoluene	ND	1.0	ug/l
Methylene chloride	ND	1.0	"
Naphthalene	ND	1.0	"
n-Propylbenzene	ND	1.0	"
Styrene	ND	1.0	"
1,1,2,2-Tetrachloroethane	ND	1.0	"
1,1,1,2-Tetrachloroethane	ND	1.0	"
Tetrachloroethene	ND	1.0	"
1,2,3-Trichlorobenzene	ND	1.0	"
1,2,4-Trichlorobenzene	ND	1.0	"
1,1,2-Trichloroethane	ND	1.0	"
1,1,1-Trichloroethane	ND	1.0	"
Trichloroethene	ND	1.0	"
Trichlorofluoromethane	ND	1.0	"
1,2,3-Trichloropropane	ND	1.0	"
1,3,5-Trimethylbenzene	ND	1.0	"
1,2,4-Trimethylbenzene	ND	1.0	"
Vinyl chloride	ND	1.0	"
Benzene	ND	0.50	"
Toluene	ND	0.50	"
Ethylbenzene	ND	0.50	"
m,p-Xylene	ND	1.0	"
o-Xylene	ND	0.50	"
Tert-amyl methyl ether	ND	2.0	"
Tert-butyl alcohol	ND	10	"
Di-isopropyl ether	ND	2.0	"
Ethyl tert-butyl ether	ND	2.0	"
Methyl tert-butyl ether	ND	1.0	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"
Surrogate 4-Bromofluorobenzene	8.52		"
Surrogate Dibromofluoromethane	8.19		"
Surrogate Toluene-d8	7.94		"

8.00 106 83.5-119
8.00 102 81-136
8.00 99.2 88.8-117

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040222 - EPA 5030 GCMS

LCS (3040222-BS1)

Prepared: 04/02/13 Analyzed: 04/03/13

Chlorobenzene	18.8	1.0	ug/l	20.0		94.0	75-125			
1,1-Dichloroethene	23.2	1.0	"	20.0		116	75-125			
Trichloroethene	19.1	1.0	"	20.0		95.6	75-125			
Benzene	20.9	0.50	"	20.0		105	75-125			
Toluene	18.4	0.50	"	20.0		92.2	75-125			
Surrogate 4-Bromofluorobenzene	8.29		"	8.00		104	83.5-119			
Surrogate Dibromofluoromethane	10.1		"	8.00		126	81-136			
Surrogate Toluene-d8	7.61		"	8.00		95.1	88.8-117			

Matrix Spike (3040222-MS1)

Source: T130750-01

Prepared: 04/02/13 Analyzed: 04/03/13

Chlorobenzene	18.5	1.0	ug/l	20.0	ND	92.6	75-125			
1,1-Dichloroethene	21.6	1.0	"	20.0	ND	108	75-125			
Trichloroethene	17.9	1.0	"	20.0	ND	89.4	75-125			
Benzene	53.8	0.50	"	20.0	34.2	98.4	75-125			
Toluene	19.6	0.50	"	20.0	1.93	88.2	75-125			
Surrogate 4-Bromofluorobenzene	8.63		"	8.00		108	83.5-119			
Surrogate Dibromofluoromethane	9.37		"	8.00		117	81-136			
Surrogate Toluene-d8	7.81		"	8.00		97.6	88.8-117			

Matrix Spike Dup (3040222-MSD1)

Source: T130750-01

Prepared: 04/02/13 Analyzed: 04/03/13

Chlorobenzene	19.6	1.0	ug/l	20.0	ND	98.0	75-125	5.77	20	
1,1-Dichloroethene	24.3	1.0	"	20.0	ND	122	75-125	11.6	20	
Trichloroethene	19.0	1.0	"	20.0	ND	95.2	75-125	6.39	20	
Benzene	64.1	0.50	"	20.0	34.2	150	75-125	17.3	20	QM-05
Toluene	21.1	0.50	"	20.0	1.93	95.8	75-125	7.48	20	
Surrogate 4-Bromofluorobenzene	8.25		"	8.00		103	83.5-119			
Surrogate Dibromofluoromethane	10.4		"	8.00		130	81-136			
Surrogate Toluene-d8	7.52		"	8.00		94.0	88.8-117			

SunStar Laboratories, Inc.

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/09/13 09:44
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Conventional Chemistry Parameters by APHA/EPA/ASTM Methods - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040231 - General Preparation

Duplicate (3040231-DUP1) **Source: T130756-01** Prepared & Analyzed: 04/02/13

Total Alkalinity	465	20	mg/l		465			0.00	25	
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Anions by EPA Method 300.0 - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040228 - General Preparation

Blank (3040228-BLK1)				Prepared & Analyzed: 04/02/13						
Nitrate as N	ND	0.100	mg/l							
Nitrite as N	ND	0.100	"							
Nitrite as NO2	ND	0.500	"							
Nitrate as NO3	ND	0.500	"							

LCS (3040228-BS1)				Prepared & Analyzed: 04/02/13						
Nitrate as NO3	0.516	0.500	mg/l	0.500		103	75-125			

Matrix Spike (3040228-MS1)				Source: T130754-04		Prepared & Analyzed: 04/02/13				
Nitrate as NO3	0.493	0.500	mg/l	0.500	ND	98.6	75-125			

Matrix Spike Dup (3040228-MSD1)				Source: T130754-04		Prepared & Analyzed: 04/02/13				
Nitrate as NO3	0.515	0.500	mg/l	0.500	ND	103	75-125	4.37	20	

Batch 3040312 - General Preparation

Blank (3040312-BLK1)				Prepared & Analyzed: 04/03/13						
Sulfate as SO4	ND	0.500	mg/l							

LCS (3040312-BS1)				Prepared & Analyzed: 04/03/13						
Sulfate as SO4	9.55	0.500	mg/l	10.0		95.5	75-125			

Matrix Spike (3040312-MS1)				Source: T130772-01		Prepared & Analyzed: 04/03/13				
Sulfate as SO4	9.38	0.500	mg/l	10.0	1.23	81.5	75-125			

Matrix Spike Dup (3040312-MSD1)				Source: T130772-01		Prepared & Analyzed: 04/03/13				
Sulfate as SO4	9.18	0.500	mg/l	10.0	1.23	79.5	75-125	2.16	20	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

RSK-175 - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040216 - EPA 3810m Headspace

Blank (3040216-BLK1)

Prepared & Analyzed: 04/02/13

Methane	ND	1.00	ug/l
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Duplicate (3040216-DUP1)

Source: T130756-01

Prepared & Analyzed: 04/02/13

Methane	ND	1.00	ug/l	ND	20
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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 09:44

Notes and Definitions

QM-05 The spike recovery was outside acceptance limits for the MS and/or MSD due to possible matrix interference. The LCS was within acceptance criteria. The data is acceptable as no negative impact on data is expected.

DET Analyte DETECTED

ND Analyte NOT DETECTED at or above the reporting limit

NR Not Reported

dry Sample results reported on a dry weight basis

RPD Relative Percent Difference

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Katherine Shields, Jr. Project Manager

Chain of Custody Record

Date: 4-1-13

Page: 1 OF 1

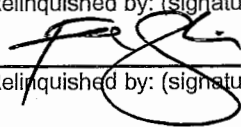
Project Name: CENCO

Collector: Frane Sosic

Client Project #: 1003-001-300

Batch #: T130756

EDF #:

Sample ID	Date Sampled	Time	Sample Type	TPHg (8015 M)	VOCs (8260 B)	Methane (8015)	Alkalinity	300 IC Nitrate, Sulfate	Ferrous Iron	Total # of containers	Comments/Preservative	Laboratory ID #
LL-14A-040113	4-1-13	1045	GW	X	X	X	X	X	X	12		0
LL-14B-040113	4-1-13	1248	GW	X	X					6		0
LL-14C-040113	4-1-13	1556	GW	X	X					6		0
LL-TB-040113			Water		X					2		0
LL-W10-040113	4-1-13	1610	GW	X	X					6		05
Relinquished by: (signature)	Date / Time	Received by: (Sign / Date / Time)			Total # of containers		FS 26 32		Notes			
	F. Sasic 4-1-2013 16:15	1610 4/1/13 16:35			Chain of Custody seals		N					
Relinquished by: (signature)	Date / Time	Received by: (Sign / Date / Time)			Seals intact? Y/N/NA		N/A					
					Received good condition/cold		Y		3.2°			
Relinquished by: (signature)	Date / Time	Received by: (Sign / Date / Time)			Turn around time:		Standard					

Sample disposal instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____

SAMPLE RECEIVING REVIEW SHEET

BATCH # T130756

Client Name: Murex

Project: Cenco

Received by: Patrick

Date/Time Received: 4/1/13 1635

Delivered by: ☐ Client ☒ SunStar Courier ☐ GSO ☐ FedEx ☐ Other _____

Total number of coolers received 1

Temp criteria = 6°C > 0°C (no frozen containers)

Temperature: cooler #1 3.4 °C +/- the CF (- 0.2°C) = 3.2 °C corrected temperature

cooler #2 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

cooler #3 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

Samples outside temp. but received on ice, w/in 6 hours of final sampling. ☒ Yes ☐ No* ☐ N/A

Custody Seals Intact on Cooler/Sample ☐ Yes ☐ No* ☒ N/A

Sample Containers Intact ☒ Yes ☐ No*

Sample labels match COC ID's ☒ Yes ☐ No*

Total number of containers received match COC ☒ Yes ☐ No*

Proper containers received for analyses requested on COC ☒ Yes ☐ No*

Proper preservative indicated on COC/containers for analyses requested ☒ Yes ☐ No* ☐ N/A

Complete shipment received in good condition with correct temperatures, containers, labels, volumes preservatives and within method specified holding times. ☒ Yes ☐ No*

* Complete Non-Conformance Receiving Sheet if checked

Cooler/Sample Review - Initials and date [Signature] 4/1/13

Comments:



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09 April 2013

Jeremy Squire
Murex
15375 Barranca Parkway, Suite K-101
Irvine, CA 92861
RE: Cenco

Enclosed are the results of analyses for samples received by the laboratory on 04/02/13 16:13. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Katherine Shields
Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
LL_15A_040213	T130772-01	Water	04/02/13 11:29	04/02/13 16:13
LL_15B_040213	T130772-02	Water	04/02/13 12:45	04/02/13 16:13
LL_15C_040213	T130772-03	Water	04/02/13 14:36	04/02/13 16:13
LL_W12_040213	T130772-04	Water	04/02/13 16:00	04/02/13 16:13
LL_TB_040213	T130772-05	Water	04/02/13 00:00	04/02/13 16:13

SunStar Laboratories, Inc.

Katherine Shields

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Katherine Shields, Jr. Project Manager



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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

LL_15A_040213
T130772-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	3400	50	ug/l	1	3040317	04/03/13	04/04/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		93.0 %	65-135		"	"	"	"	

Metals by SM 3500 Series Methods

Ferrous Iron	0.383	0.100	mg/l	1	3040247	04/02/13	04/08/13	EPA6010/S M3500	
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Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040316	04/03/13	04/09/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	4.2	1.0	"	"	"	"	"	"	
sec-Butylbenzene	2.9	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

LL_15A_040213
T130772-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1-Dichloroethene	ND	1.0	ug/l	1	3040316	04/03/13	04/09/13	EPA 8260B	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	8.5	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	2.4	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	260	1.0	"	"	"	"	"	"	E-1
n-Propylbenzene	19	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	67	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	230	1.0	"	"	"	"	"	"	E-1
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	86	0.50	"	"	"	"	"	"	
Toluene	32	0.50	"	"	"	"	"	"	
Ethylbenzene	79	0.50	"	"	"	"	"	"	
m,p-Xylene	460	25	"	25	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

LL_15A_040213
T130772-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

o-Xylene	130	0.50	ug/l	1	3040316	04/03/13	04/09/13	EPA 8260B	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	120	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	72	1.0	"	"	"	"	"	"	
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene		102 %	83.5-119		"	"	"	"	
Surrogate: Dibromofluoromethane		100 %	81-136		"	"	"	"	
Surrogate: Toluene-d8		101 %	88.8-117		"	"	"	"	

Conventional Chemistry Parameters by APHA/EPA/ASTM Methods

Total Alkalinity	790	20	mg/l	1	3040308	04/03/13	04/03/13	EPA 310.1	
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Anions by EPA Method 300.0

Sulfate as SO4	1.23	0.500	mg/l	1	3040312	04/03/13	04/03/13	EPA 300.0	
Nitrate as NO3	ND	0.500	"	"	"	"	"	"	

RSK-175

Methane	1.99	1.00	ug/l	1	3040313	04/03/13	04/04/13	RSK-175	
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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

LL_15B_040213
T130772-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	430	50	ug/l	1	3040317	04/03/13	04/04/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene	94.7 %	65-135	"	"	"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040316	04/03/13	04/09/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	1.3	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

LL_15B_040213
T130772-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,3-Dichloropropane	ND	1.0	ug/l	1	3040316	04/03/13	04/09/13	EPA 8260B	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	2.4	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	46	1.0	"	"	"	"	"	"	
n-Propylbenzene	5.8	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	19	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	62	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	1.0	0.50	"	"	"	"	"	"	
Toluene	2.3	0.50	"	"	"	"	"	"	
Ethylbenzene	13	0.50	"	"	"	"	"	"	
m,p-Xylene	87	1.0	"	"	"	"	"	"	
o-Xylene	19	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	180	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

LL_15B_040213
T130772-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Ethyl tert-butyl ether	ND	2.0	ug/l	1	3040316	04/03/13	04/09/13	EPA 8260B
Methyl tert-butyl ether	19	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		109 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		99.0 %	81-136		"	"	"	"
Surrogate: Toluene-d8		103 %	88.8-117		"	"	"	"

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

LL_15C_040213
T130772-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	410	50	ug/l	1	3040317	04/03/13	04/04/13	EPA 8015C
Surrogate: 4-Bromofluorobenzene	92.0 %	65-135	"	"	"	"	"	"

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040316	04/03/13	04/09/13	EPA 8260B
Bromochloromethane	ND	1.0	"	"	"	"	"	"
Bromodichloromethane	ND	1.0	"	"	"	"	"	"
Bromoform	ND	1.0	"	"	"	"	"	"
Bromomethane	ND	1.0	"	"	"	"	"	"
n-Butylbenzene	ND	1.0	"	"	"	"	"	"
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"
Chlorobenzene	ND	1.0	"	"	"	"	"	"
Chloroethane	ND	1.0	"	"	"	"	"	"
Chloroform	ND	1.0	"	"	"	"	"	"
Chloromethane	ND	1.0	"	"	"	"	"	"
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"
Dibromochloromethane	ND	1.0	"	"	"	"	"	"
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"
Dibromomethane	ND	1.0	"	"	"	"	"	"
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

LL_15C_040213
T130772-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040316	04/03/13	04/09/13	EPA 8260B	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	1.2	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	30	1.0	"	"	"	"	"	"	
n-Propylbenzene	2.9	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	13	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	42	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	20	0.50	"	"	"	"	"	"	
Toluene	5.8	0.50	"	"	"	"	"	"	
Ethylbenzene	9.8	0.50	"	"	"	"	"	"	
m,p-Xylene	86	1.0	"	"	"	"	"	"	
o-Xylene	21	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	25	10	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Cenco
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Reported:
04/09/13 16:50

LL_15C_040213
T130772-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Di-isopropyl ether	ND	2.0	ug/l	1	3040316	04/03/13	04/09/13	EPA 8260B
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	6.3	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		112 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		98.0 %	81-136		"	"	"	"
Surrogate: Toluene-d8		104 %	88.8-117		"	"	"	"

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

LL_W12_040213
T130772-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	150	50	ug/l	1	3040317	04/03/13	04/04/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		90.9 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040316	04/03/13	04/09/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	1.0	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

LL_W12_040213
T130772-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,3-Dichloropropane	ND	1.0	ug/l	1	3040316	04/03/13	04/09/13	EPA 8260B	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	1.2	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	13	1.0	"	"	"	"	"	"	
n-Propylbenzene	3.0	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	4.6	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	15	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	ND	0.50	"	"	"	"	"	"	
Toluene	ND	0.50	"	"	"	"	"	"	
Ethylbenzene	3.0	0.50	"	"	"	"	"	"	
m,p-Xylene	17	1.0	"	"	"	"	"	"	
o-Xylene	4.2	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	26	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	

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Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

LL_W12_040213
T130772-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Ethyl tert-butyl ether	ND	2.0	ug/l	1	3040316	04/03/13	04/09/13	EPA 8260B
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		112 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		100 %	81-136		"	"	"	"
Surrogate: Toluene-d8		102 %	88.8-117		"	"	"	"

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

LL_TB_040213
T130772-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040316	04/03/13	04/09/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

LL_TB_040213
T130772-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

cis-1,3-Dichloropropene	ND	0.50	ug/l	1	3040316	04/03/13	04/09/13	EPA 8260B	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	ND	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	ND	1.0	"	"	"	"	"	"	
n-Propylbenzene	ND	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	ND	0.50	"	"	"	"	"	"	
Toluene	ND	0.50	"	"	"	"	"	"	
Ethylbenzene	ND	0.50	"	"	"	"	"	"	
m,p-Xylene	ND	1.0	"	"	"	"	"	"	
o-Xylene	ND	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	ND	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"	

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Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

LL_TB_040213
T130772-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	ug/l	1	3040316	04/03/13	04/09/13	EPA 8260B	
Surrogate: 4-Bromofluorobenzene	110 %	83.5-119			"	"	"	"	
Surrogate: Dibromofluoromethane	100 %	81-136			"	"	"	"	
Surrogate: Toluene-d8	103 %	88.8-117			"	"	"	"	

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/09/13 16:50
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Purgeable Petroleum Hydrocarbons by EPA 8015C - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040317 - EPA 5030 GC

Blank (3040317-BLK1)

Prepared: 04/03/13 Analyzed: 04/04/13

C6-C12 (GRO)	ND	50	ug/l							
Surrogate 4-Bromofluorobenzene	88.0		"	100		88.0	65-135			

LCS (3040317-BS1)

Prepared: 04/03/13 Analyzed: 04/04/13

C6-C12 (GRO)	5400	50	ug/l	5520		97.8	75-125			
Surrogate 4-Bromofluorobenzene	68.9		"	100		68.9	65-135			

LCS Dup (3040317-BSD1)

Prepared: 04/03/13 Analyzed: 04/04/13

C6-C12 (GRO)	4830	50	ug/l	5520		87.5	75-125	11.0	20	
Surrogate 4-Bromofluorobenzene	85.4		"	100		85.4	65-135			

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

Metals by SM 3500 Series Methods - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040247 - EPA 3010A

Blank (3040247-BLK1)

Prepared: 04/02/13 Analyzed: 04/08/13

Ferrous Iron	ND	0.100	mg/l							
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Duplicate (3040247-DUP1)

Source: T130772-01

Prepared: 04/02/13 Analyzed: 04/08/13

Ferrous Iron	0.326	0.100	mg/l		0.383			16.0	200	
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SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Katherine Shields

Katherine Shields, Jr. Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040316 - EPA 5030 GCMS

Blank (3040316-BLK1)

Prepared: 04/03/13 Analyzed: 04/08/13

Bromobenzene	ND	1.0	ug/l
Bromochloromethane	ND	1.0	"
Bromodichloromethane	ND	1.0	"
Bromoform	ND	1.0	"
Bromomethane	ND	1.0	"
n-Butylbenzene	ND	1.0	"
sec-Butylbenzene	ND	1.0	"
tert-Butylbenzene	ND	1.0	"
Carbon tetrachloride	ND	0.50	"
Chlorobenzene	ND	1.0	"
Chloroethane	ND	1.0	"
Chloroform	ND	1.0	"
Chloromethane	ND	1.0	"
2-Chlorotoluene	ND	1.0	"
4-Chlorotoluene	ND	1.0	"
Dibromochloromethane	ND	1.0	"
1,2-Dibromo-3-chloropropane	ND	1.0	"
1,2-Dibromoethane (EDB)	ND	1.0	"
Dibromomethane	ND	1.0	"
1,2-Dichlorobenzene	ND	1.0	"
1,3-Dichlorobenzene	ND	1.0	"
1,4-Dichlorobenzene	ND	1.0	"
Dichlorodifluoromethane	ND	0.50	"
1,1-Dichloroethane	ND	1.0	"
1,2-Dichloroethane	ND	0.50	"
1,1-Dichloroethene	ND	1.0	"
cis-1,2-Dichloroethene	ND	1.0	"
trans-1,2-Dichloroethene	ND	1.0	"
1,2-Dichloropropane	ND	1.0	"
1,3-Dichloropropane	ND	1.0	"
2,2-Dichloropropane	ND	1.0	"
1,1-Dichloropropene	ND	1.0	"
cis-1,3-Dichloropropene	ND	0.50	"
trans-1,3-Dichloropropene	ND	0.50	"
Hexachlorobutadiene	ND	1.0	"
Isopropylbenzene	ND	1.0	"

SunStar Laboratories, Inc.

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Katherine Shields, Jr. Project Manager



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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040316 - EPA 5030 GCMS

Blank (3040316-BLK1)

Prepared: 04/03/13 Analyzed: 04/08/13

p-Isopropyltoluene	ND	1.0	ug/l
Methylene chloride	ND	1.0	"
Naphthalene	ND	1.0	"
n-Propylbenzene	ND	1.0	"
Styrene	ND	1.0	"
1,1,2,2-Tetrachloroethane	ND	1.0	"
1,1,1,2-Tetrachloroethane	ND	1.0	"
Tetrachloroethene	ND	1.0	"
1,2,3-Trichlorobenzene	ND	1.0	"
1,2,4-Trichlorobenzene	ND	1.0	"
1,1,2-Trichloroethane	ND	1.0	"
1,1,1-Trichloroethane	ND	1.0	"
Trichloroethene	ND	1.0	"
Trichlorofluoromethane	ND	1.0	"
1,2,3-Trichloropropane	ND	1.0	"
1,3,5-Trimethylbenzene	ND	1.0	"
1,2,4-Trimethylbenzene	ND	1.0	"
Vinyl chloride	ND	1.0	"
Benzene	ND	0.50	"
Toluene	ND	0.50	"
Ethylbenzene	ND	0.50	"
m,p-Xylene	ND	1.0	"
o-Xylene	ND	0.50	"
Tert-amyl methyl ether	ND	2.0	"
Tert-butyl alcohol	ND	10	"
Di-isopropyl ether	ND	2.0	"
Ethyl tert-butyl ether	ND	2.0	"
Methyl tert-butyl ether	ND	1.0	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"
Surrogate 4-Bromofluorobenzene	8.70		"
Surrogate Dibromofluoromethane	7.86		"
Surrogate Toluene-d8	8.37		"

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040316 - EPA 5030 GCMS

LCS (3040316-BS1)

Prepared: 04/03/13 Analyzed: 04/09/13

Chlorobenzene	19.2	1.0	ug/l	20.0		95.8	75-125			
1,1-Dichloroethene	20.5	1.0	"	20.0		103	75-125			
Trichloroethene	18.4	1.0	"	20.0		92.2	75-125			
Benzene	19.9	0.50	"	20.0		99.6	75-125			
Toluene	18.0	0.50	"	20.0		89.8	75-125			
Surrogate 4-Bromofluorobenzene	8.40		"	8.00		105	83.5-119			
Surrogate Dibromofluoromethane	8.53		"	8.00		107	81-136			
Surrogate Toluene-d8	8.28		"	8.00		104	88.8-117			

LCS Dup (3040316-BSD1)

Prepared: 04/03/13 Analyzed: 04/09/13

Chlorobenzene	18.6	1.0	ug/l	20.0		93.0	75-125	3.07	20	
1,1-Dichloroethene	19.6	1.0	"	20.0		98.0	75-125	4.63	20	
Trichloroethene	18.0	1.0	"	20.0		90.0	75-125	2.42	20	
Benzene	18.8	0.50	"	20.0		94.2	75-125	5.52	20	
Toluene	17.4	0.50	"	20.0		86.9	75-125	3.28	20	
Surrogate 4-Bromofluorobenzene	8.44		"	8.00		106	83.5-119			
Surrogate Dibromofluoromethane	8.48		"	8.00		106	81-136			
Surrogate Toluene-d8	8.07		"	8.00		101	88.8-117			

SunStar Laboratories, Inc.

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/09/13 16:50
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Conventional Chemistry Parameters by APHA/EPA/ASTM Methods - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040308 - General Preparation

Duplicate (3040308-DUP1) **Source: T130772-01** Prepared & Analyzed: 04/03/13

Total Alkalinity	800	20	mg/l		790			1.26	25	
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SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

Anions by EPA Method 300.0 - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040312 - General Preparation

Blank (3040312-BLK1)

Prepared & Analyzed: 04/03/13

Sulfate as SO ₄	ND	0.500	mg/l
Nitrate as NO ₃	ND	0.500	"

LCS (3040312-BS1)

Prepared & Analyzed: 04/03/13

Sulfate as SO ₄	9.55	0.500	mg/l	10.0	95.5	75-125
Nitrate as NO ₃	0.521	0.500	"	0.500	104	75-125

Matrix Spike (3040312-MS1)

Source: T130772-01

Prepared & Analyzed: 04/03/13

Sulfate as SO ₄	9.38	0.500	mg/l	10.0	1.23	81.5	75-125
Nitrate as NO ₃	0.532	0.500	"	0.500	ND	106	75-125

Matrix Spike Dup (3040312-MSD1)

Source: T130772-01

Prepared & Analyzed: 04/03/13

Sulfate as SO ₄	9.18	0.500	mg/l	10.0	1.23	79.5	75-125	2.16	20
Nitrate as NO ₃	0.521	0.500	"	0.500	ND	104	75-125	2.09	20

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/09/13 16:50

RSK-175 - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040313 - EPA 3810m Headspace

Blank (3040313-BLK1)

Prepared: 04/03/13 Analyzed: 04/04/13

Methane	ND	1.00	ug/l							
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Duplicate (3040313-DUP1)

Source: T130772-01

Prepared: 04/03/13 Analyzed: 04/04/13

Methane	1.90	1.00	ug/l		1.99			4.63	20	
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SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/09/13 16:50
--	--	------------------------------------

Notes and Definitions

E-1	The final dilution was lower than the original data or previous dilutions. The highest recovered concentration was reported even though it was above calibration range.
DET	Analyte DETECTED
ND	Analyte NOT DETECTED at or above the reporting limit
NR	Not Reported
dry	Sample results reported on a dry weight basis
RPD	Relative Percent Difference

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager

SunStar Laboratories, Inc.
25712 Commercentre Dr
Lake Forest, CA 92630
949-297-5020

Chain of Custody Record

Client: MUREX ENVIRONMENTAL INC.

Address: 2640 Walnut Ave, Unit F

Phone: (714) 508-0800 Fax: (714) 508-0880

Project Manager: Jeremy Squire (714) 604-5836

Date:

Project Name: CENCO

Collector: Frane Sosic

Batch #: T130772

Page:

OF

Client Project #: 1003-001-300

EDF #:

[illegible]

Sample disposal instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____

SAMPLE RECEIVING REVIEW SHEET

BATCH # 7130772

Client Name: Murex

Project: Cenco

Received by: Patrick

Date/Time Received: 4/2/13 1613

Delivered by: ☐ Client ☒ SunStar Courier ☐ GSO ☐ FedEx ☐ Other _____

Total number of coolers received 1 Temp criteria = 6°C > 0°C (no frozen containers)

Temperature: cooler #1 3.0 °C +/- the CF (-0.2°C) = 2.8 °C corrected temperature

cooler #2 _____ °C +/- the CF (-0.2°C) = _____ °C corrected temperature

cooler #3 _____ °C +/- the CF (-0.2°C) = _____ °C corrected temperature

Samples outside temp. but received on ice, w/in 6 hours of final sampling. ☒ Yes ☐ No* ☐ N/A

Custody Seals Intact on Cooler/Sample ☐ Yes ☐ No* ☒ N/A

Sample Containers Intact ☒ Yes ☐ No*

Sample labels match COC ID's ☒ Yes ☐ No*

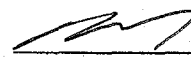
Total number of containers received match COC ☒ Yes ☐ No*

Proper containers received for analyses requested on COC ☒ Yes ☐ No*

Proper preservative indicated on COC/containers for analyses requested ☒ Yes ☐ No* ☐ N/A

Complete shipment received in good condition with correct temperatures, containers, labels, volumes preservatives and within method specified holding times. ☐ Yes ☐ No*

* Complete Non-Conformance Receiving Sheet if checked

Cooler/Sample Review - Initials and date  4/2/13

Comments:



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Lake Forest, California 92630
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949.297.5027 Fax

10 April 2013

Jeremy Squire
Murex
15375 Barranca Parkway, Suite K-101
Irvine, CA 92861
RE: Cenco

Enclosed are the results of analyses for samples received by the laboratory on 04/03/13 16:30. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Wendy Hsiao For Katherine Shields
Jr. Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
LL_16A_040313	T130785-01	Water	04/03/13 10:47	04/03/13 16:30
LL_16B_040313	T130785-02	Water	04/03/13 12:03	04/03/13 16:30
LL_16C_040313	T130785-03	Water	04/03/13 14:35	04/03/13 16:30
LL_W1_040313	T130785-04	Water	04/03/13 15:48	04/03/13 16:30
LL_TB_040313	T130785-05	Water	04/03/13 00:00	04/03/13 16:30

SunStar Laboratories, Inc.

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Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

LL_16A_040313
T130785-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	340	50	ug/l	1	3040410	04/04/13	04/08/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		128 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040409	04/04/13	04/08/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	1.8	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	1.2	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	2.0	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Wendy Hsiao

Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

LL_16A_040313
T130785-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040409	04/04/13	04/08/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	5.5	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	20	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

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Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex
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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

LL_16A_040313
T130785-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3040409	04/04/13	04/08/13	EPA 8260B	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"	
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene		127 %	83.5-119		"	"	"	"	S-GC
Surrogate: Dibromofluoromethane		101 %	81-136		"	"	"	"	
Surrogate: Toluene-d8		99.9 %	88.8-117		"	"	"	"	

SunStar Laboratories, Inc.

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Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

LL_16B_040313
T130785-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	ND	50	ug/l	1	3040410	04/04/13	04/08/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		124 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040409	04/04/13	04/08/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	1.9	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	1.2	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

LL_16B_040313
T130785-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,3-Dichloropropane	ND	1.0	ug/l	1	3040409	04/04/13	04/08/13	EPA 8260B
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	9.1	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	1.5	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	3.4	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	1.1	1.0	"	"	"	"	"	"
o-Xylene	0.79	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"
Tert-butyl alcohol	ND	10	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

LL_16B_040313
T130785-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Di-isopropyl ether	ND	2.0	ug/l	1	3040409	04/04/13	04/08/13	EPA 8260B
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		116 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		99.1 %	81-136		"	"	"	"
Surrogate: Toluene-d8		97.5 %	88.8-117		"	"	"	"

SunStar Laboratories, Inc.

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/10/13 16:29
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LL_16C_040313
T130785-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	ND	50	ug/l	1	3040410	04/04/13	04/08/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		125 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040409	04/04/13	04/08/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	2.6	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

LL_16C_040313
T130785-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040409	04/04/13	04/08/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	7.8	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	0.70	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	0.56	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

LL_16C_040313
T130785-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3040409	04/04/13	04/08/13	EPA 8260B
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		111 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		106 %	81-136		"	"	"	"
Surrogate: Toluene-d8		97.6 %	88.8-117		"	"	"	"

RSK-175

Methane	6.22	1.00	ug/l	1	3040417	04/04/13	04/04/13	RSK-175
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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

LL_W1_040313
T130785-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	ND	50	ug/l	1	3040410	04/04/13	04/08/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		136 %	65-135		"	"	"	"	S-04

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040409	04/04/13	04/09/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

LL_W1_040313
T130785-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040409	04/04/13	04/09/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	5.6	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	1.2	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

LL_W1_040313
T130785-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3040409	04/04/13	04/09/13	EPA 8260B
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		107 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		102 %	81-136		"	"	"	"
Surrogate: Toluene-d8		96.9 %	88.8-117		"	"	"	"

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

LL_TB_040313
T130785-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040409	04/04/13	04/09/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

LL_TB_040313
T130785-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

cis-1,3-Dichloropropene	ND	0.50	ug/l	1	3040409	04/04/13	04/09/13	EPA 8260B	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	ND	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	ND	1.0	"	"	"	"	"	"	
n-Propylbenzene	ND	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	ND	0.50	"	"	"	"	"	"	
Toluene	ND	0.50	"	"	"	"	"	"	
Ethylbenzene	ND	0.50	"	"	"	"	"	"	
m,p-Xylene	ND	1.0	"	"	"	"	"	"	
o-Xylene	ND	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	ND	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

LL_TB_040313
T130785-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	ug/l	1	3040409	04/04/13	04/09/13	EPA 8260B	
Surrogate: 4-Bromofluorobenzene	109 %	83.5-119			"	"	"	"	
Surrogate: Dibromofluoromethane	100 %	81-136			"	"	"	"	
Surrogate: Toluene-d8	98.4 %	88.8-117			"	"	"	"	

SunStar Laboratories, Inc.

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/10/13 16:29
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Purgeable Petroleum Hydrocarbons by EPA 8015C - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
Batch 3040410 - EPA 5030 GC										
Blank (3040410-BLK1)				Prepared: 04/04/13 Analyzed: 04/08/13						
C6-C12 (GRO)	ND	50	ug/l							
Surrogate 4-Bromofluorobenzene	121		"	100		121	65-135			
LCS (3040410-BS1)				Prepared: 04/04/13 Analyzed: 04/08/13						
C6-C12 (GRO)	5350	50	ug/l	5520		97.0	75-125			
Surrogate 4-Bromofluorobenzene	101		"	100		101	65-135			
Matrix Spike (3040410-MS1)				Source: T130785-01		Prepared: 04/04/13 Analyzed: 04/08/13				
C6-C12 (GRO)	4030	50	ug/l	5520	343	66.8	65-135			
Surrogate 4-Bromofluorobenzene	88.1		"	100		88.1	65-135			
Matrix Spike Dup (3040410-MSD1)				Source: T130785-01		Prepared: 04/04/13 Analyzed: 04/08/13				
C6-C12 (GRO)	3640	50	ug/l	5520	343	59.8	65-135	10.0	20	QM-05
Surrogate 4-Bromofluorobenzene	89.0		"	100		89.0	65-135			

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040409 - EPA 5030 GCMS

Blank (3040409-BLK1)

Prepared: 04/04/13 Analyzed: 04/08/13

Bromobenzene	ND	1.0	ug/l
Bromochloromethane	ND	1.0	"
Bromodichloromethane	ND	1.0	"
Bromoform	ND	1.0	"
Bromomethane	ND	1.0	"
n-Butylbenzene	ND	1.0	"
sec-Butylbenzene	ND	1.0	"
tert-Butylbenzene	ND	1.0	"
Carbon tetrachloride	ND	0.50	"
Chlorobenzene	ND	1.0	"
Chloroethane	ND	1.0	"
Chloroform	ND	1.0	"
Chloromethane	ND	1.0	"
2-Chlorotoluene	ND	1.0	"
4-Chlorotoluene	ND	1.0	"
Dibromochloromethane	ND	1.0	"
1,2-Dibromo-3-chloropropane	ND	1.0	"
1,2-Dibromoethane (EDB)	ND	1.0	"
Dibromomethane	ND	1.0	"
1,2-Dichlorobenzene	ND	1.0	"
1,3-Dichlorobenzene	ND	1.0	"
1,4-Dichlorobenzene	ND	1.0	"
Dichlorodifluoromethane	ND	0.50	"
1,1-Dichloroethane	ND	1.0	"
1,2-Dichloroethane	ND	0.50	"
1,1-Dichloroethene	ND	1.0	"
cis-1,2-Dichloroethene	ND	1.0	"
trans-1,2-Dichloroethene	ND	1.0	"
1,2-Dichloropropane	ND	1.0	"
1,3-Dichloropropane	ND	1.0	"
2,2-Dichloropropane	ND	1.0	"
1,1-Dichloropropene	ND	1.0	"
cis-1,3-Dichloropropene	ND	0.50	"
trans-1,3-Dichloropropene	ND	0.50	"
Hexachlorobutadiene	ND	1.0	"
Isopropylbenzene	ND	1.0	"

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040409 - EPA 5030 GCMS

Blank (3040409-BLK1)

Prepared: 04/04/13 Analyzed: 04/08/13

p-Isopropyltoluene	ND	1.0	ug/l
Methylene chloride	ND	1.0	"
Naphthalene	ND	1.0	"
n-Propylbenzene	ND	1.0	"
Styrene	ND	1.0	"
1,1,2,2-Tetrachloroethane	ND	1.0	"
1,1,1,2-Tetrachloroethane	ND	1.0	"
Tetrachloroethene	ND	1.0	"
1,2,3-Trichlorobenzene	ND	1.0	"
1,2,4-Trichlorobenzene	ND	1.0	"
1,1,2-Trichloroethane	ND	1.0	"
1,1,1-Trichloroethane	ND	1.0	"
Trichloroethene	ND	1.0	"
Trichlorofluoromethane	ND	1.0	"
1,2,3-Trichloropropane	ND	1.0	"
1,3,5-Trimethylbenzene	ND	1.0	"
1,2,4-Trimethylbenzene	ND	1.0	"
Vinyl chloride	ND	1.0	"
Benzene	ND	0.50	"
Toluene	ND	0.50	"
Ethylbenzene	ND	0.50	"
m,p-Xylene	ND	1.0	"
o-Xylene	ND	0.50	"
Tert-amyl methyl ether	ND	2.0	"
Tert-butyl alcohol	ND	10	"
Di-isopropyl ether	ND	2.0	"
Ethyl tert-butyl ether	ND	2.0	"
Methyl tert-butyl ether	ND	1.0	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"
Surrogate 4-Bromofluorobenzene	8.92		"
Surrogate Dibromofluoromethane	8.12		"
Surrogate Toluene-d8	7.54		"

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040409 - EPA 5030 GCMS

LCS (3040409-BS1)

Prepared: 04/04/13 Analyzed: 04/09/13

Chlorobenzene	18.8	1.0	ug/l	20.0		93.8	75-125			
1,1-Dichloroethene	24.7	1.0	"	20.0		124	75-125			
Trichloroethene	21.5	1.0	"	20.0		107	75-125			
Benzene	21.6	0.50	"	20.0		108	75-125			
Toluene	19.2	0.50	"	20.0		96.1	75-125			
Surrogate 4-Bromofluorobenzene	8.16		"	8.00		102	83.5-119			
Surrogate Dibromofluoromethane	8.37		"	8.00		105	81-136			
Surrogate Toluene-d8	7.55		"	8.00		94.4	88.8-117			

Matrix Spike (3040409-MS1)

Source: T130785-01

Prepared: 04/04/13 Analyzed: 04/09/13

Chlorobenzene	19.6	1.0	ug/l	20.0	ND	97.8	75-125			
1,1-Dichloroethene	25.7	1.0	"	20.0	ND	129	75-125			QM-05
Trichloroethene	21.6	1.0	"	20.0	ND	108	75-125			
Benzene	40.8	0.50	"	20.0	20.1	103	75-125			
Toluene	21.2	0.50	"	20.0	ND	106	75-125			
Surrogate 4-Bromofluorobenzene	8.52		"	8.00		106	83.5-119			
Surrogate Dibromofluoromethane	8.48		"	8.00		106	81-136			
Surrogate Toluene-d8	7.89		"	8.00		98.6	88.8-117			

Matrix Spike Dup (3040409-MSD1)

Source: T130785-01

Prepared: 04/04/13 Analyzed: 04/09/13

Chlorobenzene	20.5	1.0	ug/l	20.0	ND	102	75-125	4.70	20	
1,1-Dichloroethene	26.2	1.0	"	20.0	ND	131	75-125	1.89	20	QM-05
Trichloroethene	21.3	1.0	"	20.0	ND	106	75-125	1.58	20	
Benzene	39.8	0.50	"	20.0	20.1	98.6	75-125	2.31	20	
Toluene	20.3	0.50	"	20.0	ND	102	75-125	4.48	20	
Surrogate 4-Bromofluorobenzene	8.66		"	8.00		108	83.5-119			
Surrogate Dibromofluoromethane	8.78		"	8.00		110	81-136			
Surrogate Toluene-d8	7.70		"	8.00		96.2	88.8-117			

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:29

RSK-175 - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040417 - EPA 3810m Headspace

Blank (3040417-BLK1)

Prepared & Analyzed: 04/04/13

Methane	ND	1.00	ug/l
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Duplicate (3040417-DUP1)

Source: T130785-03

Prepared & Analyzed: 04/04/13

Methane	6.64	1.00	ug/l	6.22	6.53	20
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Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/10/13 16:29
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Notes and Definitions

S-GC Surrogate recovery outside of established control limits. The data was accepted based on valid recovery of the remaining surrogate(s).

S-04 The surrogate recovery for this sample is outside of established control limits due to a sample matrix effect.

QM-05 The spike recovery was outside acceptance limits for the MS and/or MSD due to possible matrix interference. The LCS was within acceptance criteria. The data is acceptable as no negative impact on data is expected.

DET Analyte DETECTED

ND Analyte NOT DETECTED at or above the reporting limit

NR Not Reported

dry Sample results reported on a dry weight basis

RPD Relative Percent Difference

SunStar Laboratories, Inc.

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Wendy Hsiao For Katherine Shields, Jr. Project Manager

Chain of Custody Record

Date: 4-3-13
Project Name: CENCO
Collector: Frane Sosic
Batch #: 7130785

Page: 1 OF 1

Client Project #: 1003-001-300

EDF #:

Sample ID	Date Sampled	Time	Sample Type	TPHg (8015 M)	VOCs (8260 B)	Methane (8015)	Alkalinity	Nitrate, Sulfate (3010)	Ferrous Iron									Total # of containers	Comments/Preservative	Laboratory ID #
LL-16A-040313	4.3.13	1047	GW	X	X													6		01
LL-16B-040313	4.3.13	1203	GW	X	X													6		02
LL-16C-040313	4.3.13	1435	GW	X	X	X												9		03
LL-W1-040313	4.3.13	1548	GW	X	X													6		04
LL-TB-040313			Water		X													2		05
Relinquished by: (signature) <i>F. Sisk</i>	Date / Time 4.3.2013 16:30	Received by: (Sign / Date / Time) <i>P. H. P.</i> 4/13/13 16:30		Total # of containers		FSEB 29		Notes												
Relinquished by: (signature) <i>[Signature]</i>	Date / Time	Received by: (Sign / Date / Time)		Chain of Custody seals																
Relinquished by: (signature)	Date / Time	Received by: (Sign / Date / Time)		Seals intact? Y/N/NA																
				Received good condition/cold		4.7														
Relinquished by: (signature)	Date / Time	Received by: (Sign / Date / Time)		Turn around time:		Standard														

Sample disposal Instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____

SAMPLE RECEIVING REVIEW SHEET

BATCH # T130785

Client Name: MUREX

Project: CENCO

Received by: PATRICK

Date/Time Received: 4.3.13 / 16:30

Delivered by : ☐ Client ☒ SunStar Courier ☐ GSO ☐ FedEx ☐ Other _____

Total number of coolers received 0

Temp criteria = 6°C > 0°C (no frozen containers)

Temperature: cooler #1 4.9 °C +/- the CF (- 0.2°C) = 4.7 °C corrected temperature

cooler #2 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

cooler #3 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

Samples outside temp. but received on ice, w/in 6 hours of final sampling. ☒ Yes ☐ No* ☐ N/A

Custody Seals Intact on Cooler/Sample ☐ Yes ☐ No* ☒ N/A

Sample Containers Intact ☒ Yes ☐ No*

Sample labels match COC ID's ☒ Yes ☐ No*

Total number of containers received match COC ☒ Yes ☐ No*

Proper containers received for analyses requested on COC ☒ Yes ☐ No*

Proper preservative indicated on COC/containers for analyses requested ☒ Yes ☐ No* ☐ N/A

Complete shipment received in good condition with correct temperatures, containers, labels, volumes preservatives and within method specified holding times. ☒ Yes ☐ No*

* Complete Non-Conformance Receiving Sheet if checked

Cooler/Sample Review - Initials and date 82 4.3.13

Comments:



25712 Commercentre Drive
Lake Forest, California 92630
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08 April 2013

Jeremy Squire
Murex
15375 Barranca Parkway, Suite K-101
Irvine, CA 92861
RE: Cenco

Enclosed are the results of analyses for samples received by the laboratory on 04/04/13 16:26. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Katherine Shields
Jr. Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/08/13 16:19

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
LL_W7_040413	T130794-01	Water	04/04/13 11:13	04/04/13 16:26
LL_W8_040413	T130794-02	Water	04/04/13 11:43	04/04/13 16:26
LL_106A_040413	T130794-03	Water	04/04/13 12:34	04/04/13 16:26
LL_107A_040413	T130794-04	Water	04/04/13 12:34	04/04/13 16:26
LL_TB_040413	T130794-05	Water	04/04/13 00:00	04/04/13 16:26

SunStar Laboratories, Inc.

Katherine Shields

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Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/08/13 16:19

LL_W7_040413
T130794-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	ND	50	ug/l	1	3040808	04/08/13	04/08/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		118 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040510	04/05/13	04/06/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/08/13 16:19

LL_W7_040413
T130794-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040510	04/05/13	04/06/13	EPA 8260B	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	ND	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	ND	1.0	"	"	"	"	"	"	
n-Propylbenzene	ND	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	ND	0.50	"	"	"	"	"	"	
Toluene	ND	0.50	"	"	"	"	"	"	
Ethylbenzene	ND	0.50	"	"	"	"	"	"	
m,p-Xylene	ND	1.0	"	"	"	"	"	"	
o-Xylene	ND	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/08/13 16:19

LL_W7_040413
T130794-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3040510	04/05/13	04/06/13	EPA 8260B
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		108 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		118 %	81-136		"	"	"	"
Surrogate: Toluene-d8		101 %	88.8-117		"	"	"	"

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/08/13 16:19

LL_W8_040413
T130794-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	ND	50	ug/l	1	3040808	04/08/13	04/08/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		113 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040510	04/05/13	04/06/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager

Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/08/13 16:19

LL_W8_040413
T130794-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040510	04/05/13	04/06/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/08/13 16:19

LL_W8_040413
T130794-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3040510	04/05/13	04/06/13	EPA 8260B
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		108 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		114 %	81-136		"	"	"	"
Surrogate: Toluene-d8		104 %	88.8-117		"	"	"	"

SunStar Laboratories, Inc.

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Katherine Shields, Jr. Project Manager



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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/08/13 16:19

LL_106A_040413
T130794-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	480	50	ug/l	1	3040808	04/08/13	04/08/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		106 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040510	04/05/13	04/06/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	4.8	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	1.9	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/08/13 16:19

LL_106A_040413
T130794-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,3-Dichloropropane	ND	1.0	ug/l	1	3040510	04/05/13	04/06/13	EPA 8260B	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	25	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	ND	1.0	"	"	"	"	"	"	
n-Propylbenzene	17	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
Vinyl chloride	14	1.0	"	"	"	"	"	"	
Benzene	6.9	0.50	"	"	"	"	"	"	
Toluene	ND	0.50	"	"	"	"	"	"	
Ethylbenzene	ND	0.50	"	"	"	"	"	"	
m,p-Xylene	ND	1.0	"	"	"	"	"	"	
o-Xylene	ND	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	ND	10	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/08/13 16:19

LL_106A_040413
T130794-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Di-isopropyl ether	ND	2.0	ug/l	1	3040510	04/05/13	04/06/13	EPA 8260B
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		112 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		112 %	81-136		"	"	"	"
Surrogate: Toluene-d8		104 %	88.8-117		"	"	"	"

SunStar Laboratories, Inc.

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/08/13 16:19
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LL_107A_040413
T130794-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	180	50	ug/l	1	3040808	04/08/13	04/08/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene	133 %	65-135	"	"	"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040510	04/05/13	04/06/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/08/13 16:19

LL_107A_040413
T130794-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040510	04/05/13	04/06/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	71	1.0	"	"	"	"	"	"
n-Propylbenzene	1.1	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	3.2	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	15	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	2.1	0.50	"	"	"	"	"	"
Ethylbenzene	1.8	0.50	"	"	"	"	"	"
m,p-Xylene	9.6	1.0	"	"	"	"	"	"
o-Xylene	5.3	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager

Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/08/13 16:19

LL_107A_040413
T130794-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3040510	04/05/13	04/06/13	EPA 8260B
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
<i>Surrogate: 4-Bromofluorobenzene</i>		<i>112 %</i>	<i>83.5-119</i>		<i>"</i>	<i>"</i>	<i>"</i>	<i>"</i>
<i>Surrogate: Dibromofluoromethane</i>		<i>112 %</i>	<i>81-136</i>		<i>"</i>	<i>"</i>	<i>"</i>	<i>"</i>
<i>Surrogate: Toluene-d8</i>		<i>96.6 %</i>	<i>88.8-117</i>		<i>"</i>	<i>"</i>	<i>"</i>	<i>"</i>

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/08/13 16:19

LL_TB_040413
T130794-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040510	04/05/13	04/06/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/08/13 16:19

LL_TB_040413
T130794-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

cis-1,3-Dichloropropene	ND	0.50	ug/l	1	3040510	04/05/13	04/06/13	EPA 8260B	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	ND	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	ND	1.0	"	"	"	"	"	"	
n-Propylbenzene	ND	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	ND	0.50	"	"	"	"	"	"	
Toluene	ND	0.50	"	"	"	"	"	"	
Ethylbenzene	ND	0.50	"	"	"	"	"	"	
m,p-Xylene	ND	1.0	"	"	"	"	"	"	
o-Xylene	ND	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	ND	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/08/13 16:19

LL_TB_040413
T130794-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	ug/l	1	3040510	04/05/13	04/06/13	EPA 8260B	
Surrogate: 4-Bromofluorobenzene	107 %	83.5-119			"	"	"	"	
Surrogate: Dibromofluoromethane	114 %	81-136			"	"	"	"	
Surrogate: Toluene-d8	97.9 %	88.8-117			"	"	"	"	

SunStar Laboratories, Inc.

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Murex	Project: Cenco	Reported:
15375 Barranca Parkway, Suite K-101	Project Number: 1003-001-300	04/08/13 16:19
Irvine CA, 92861	Project Manager: Jeremy Squire	

Purgeable Petroleum Hydrocarbons by EPA 8015C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040808 - EPA 5030 GC

Blank (3040808-BLK1)

Prepared & Analyzed: 04/08/13

C6-C12 (GRO)	ND	50	ug/l							
Surrogate 4-Bromofluorobenzene	98.3		"	100		98.3	65-135			

LCS (3040808-BS1)

Prepared & Analyzed: 04/08/13

C6-C12 (GRO)	5070	50	ug/l	5520		91.9	75-125			
Surrogate 4-Bromofluorobenzene	82.2		"	100		82.2	65-135			

LCS Dup (3040808-BSD1)

Prepared & Analyzed: 04/08/13

C6-C12 (GRO)	5190	50	ug/l	5520		94.1	75-125	2.36	20	
Surrogate 4-Bromofluorobenzene	76.2		"	100		76.2	65-135			

SunStar Laboratories, Inc.

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Murex	Project: Cenco	
15375 Barranca Parkway, Suite K-101	Project Number: 1003-001-300	
Irvine CA, 92861	Project Manager: Jeremy Squire	Reported: 04/08/13 16:19

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040510 - EPA 5030 GCMS

Blank (3040510-BLK1)

Prepared: 04/05/13 Analyzed: 04/06/13

Bromobenzene	ND	1.0	ug/l
Bromochloromethane	ND	1.0	"
Bromodichloromethane	ND	1.0	"
Bromoform	ND	1.0	"
Bromomethane	ND	1.0	"
n-Butylbenzene	ND	1.0	"
sec-Butylbenzene	ND	1.0	"
tert-Butylbenzene	ND	1.0	"
Carbon tetrachloride	ND	0.50	"
Chlorobenzene	ND	1.0	"
Chloroethane	ND	1.0	"
Chloroform	ND	1.0	"
Chloromethane	ND	1.0	"
2-Chlorotoluene	ND	1.0	"
4-Chlorotoluene	ND	1.0	"
Dibromochloromethane	ND	1.0	"
1,2-Dibromo-3-chloropropane	ND	1.0	"
1,2-Dibromoethane (EDB)	ND	1.0	"
Dibromomethane	ND	1.0	"
1,2-Dichlorobenzene	ND	1.0	"
1,3-Dichlorobenzene	ND	1.0	"
1,4-Dichlorobenzene	ND	1.0	"
Dichlorodifluoromethane	ND	0.50	"
1,1-Dichloroethane	ND	1.0	"
1,2-Dichloroethane	ND	0.50	"
1,1-Dichloroethene	ND	1.0	"
cis-1,2-Dichloroethene	ND	1.0	"
trans-1,2-Dichloroethene	ND	1.0	"
1,2-Dichloropropane	ND	1.0	"
1,3-Dichloropropane	ND	1.0	"
2,2-Dichloropropane	ND	1.0	"
1,1-Dichloropropene	ND	1.0	"
cis-1,3-Dichloropropene	ND	0.50	"
trans-1,3-Dichloropropene	ND	0.50	"
Hexachlorobutadiene	ND	1.0	"
Isopropylbenzene	ND	1.0	"

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



25712 Commercentre Drive
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949.297.5027 Fax

Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/08/13 16:19

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040510 - EPA 5030 GCMS

Blank (3040510-BLK1)

Prepared: 04/05/13 Analyzed: 04/06/13

p-Isopropyltoluene	ND	1.0	ug/l
Methylene chloride	ND	1.0	"
Naphthalene	ND	1.0	"
n-Propylbenzene	ND	1.0	"
Styrene	ND	1.0	"
1,1,2,2-Tetrachloroethane	ND	1.0	"
1,1,1,2-Tetrachloroethane	ND	1.0	"
Tetrachloroethene	ND	1.0	"
1,2,3-Trichlorobenzene	ND	1.0	"
1,2,4-Trichlorobenzene	ND	1.0	"
1,1,2-Trichloroethane	ND	1.0	"
1,1,1-Trichloroethane	ND	1.0	"
Trichloroethene	ND	1.0	"
Trichlorofluoromethane	ND	1.0	"
1,2,3-Trichloropropane	ND	1.0	"
1,3,5-Trimethylbenzene	ND	1.0	"
1,2,4-Trimethylbenzene	ND	1.0	"
Vinyl chloride	ND	1.0	"
Benzene	ND	0.50	"
Toluene	ND	0.50	"
Ethylbenzene	ND	0.50	"
m,p-Xylene	ND	1.0	"
o-Xylene	ND	0.50	"
Tert-amyl methyl ether	ND	2.0	"
Tert-butyl alcohol	ND	10	"
Di-isopropyl ether	ND	2.0	"
Ethyl tert-butyl ether	ND	2.0	"
Methyl tert-butyl ether	ND	1.0	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"
Surrogate 4-Bromofluorobenzene	8.67		"
Surrogate Dibromofluoromethane	9.52		"
Surrogate Toluene-d8	8.01		"

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex	Project: Cenco	Reported:
15375 Barranca Parkway, Suite K-101	Project Number: 1003-001-300	04/08/13 16:19
Irvine CA, 92861	Project Manager: Jeremy Squire	

Volatil Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040510 - EPA 5030 GCMS

LCS (3040510-BS1)				Prepared: 04/05/13		Analyzed: 04/06/13				
Chlorobenzene	16.7	1.0	ug/l	20.0		83.4	75-125			
1,1-Dichloroethene	22.1	1.0	"	20.0		110	75-125			
Trichloroethene	15.4	1.0	"	20.0		76.8	75-125			
Benzene	21.7	0.50	"	20.0		109	75-125			
Toluene	20.6	0.50	"	20.0		103	75-125			
Surrogate 4-Bromofluorobenzene	7.67		"	8.00		95.9	83.5-119			
Surrogate Dibromofluoromethane	9.77		"	8.00		122	81-136			
Surrogate Toluene-d8	8.09		"	8.00		101	88.8-117			

Matrix Spike (3040510-MS1)				Source: T130794-01		Prepared: 04/05/13		Analyzed: 04/06/13		
Chlorobenzene	18.0	1.0	ug/l	20.0	ND	89.8	75-125			
1,1-Dichloroethene	22.1	1.0	"	20.0	ND	111	75-125			
Trichloroethene	13.7	1.0	"	20.0	ND	68.6	75-125			QM-05
Benzene	22.3	0.50	"	20.0	ND	112	75-125			
Toluene	19.8	0.50	"	20.0	ND	98.8	75-125			
Surrogate 4-Bromofluorobenzene	8.27		"	8.00		103	83.5-119			
Surrogate Dibromofluoromethane	10.3		"	8.00		129	81-136			
Surrogate Toluene-d8	8.22		"	8.00		103	88.8-117			

Matrix Spike Dup (3040510-MSD1)				Source: T130794-01		Prepared: 04/05/13		Analyzed: 04/06/13		
Chlorobenzene	18.3	1.0	ug/l	20.0	ND	91.4	75-125	1.66	20	
1,1-Dichloroethene	23.1	1.0	"	20.0	ND	115	75-125	4.25	20	
Trichloroethene	13.8	1.0	"	20.0	ND	69.0	75-125	0.654	20	QM-05
Benzene	23.1	0.50	"	20.0	ND	115	75-125	3.44	20	
Toluene	20.6	0.50	"	20.0	ND	103	75-125	4.26	20	
Surrogate 4-Bromofluorobenzene	8.04		"	8.00		100	83.5-119			
Surrogate Dibromofluoromethane	10.4		"	8.00		129	81-136			
Surrogate Toluene-d8	7.95		"	8.00		99.4	88.8-117			

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/08/13 16:19
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Notes and Definitions

QM-05	The spike recovery was outside acceptance limits for the MS and/or MSD due to possible matrix interference. The LCS was within acceptance criteria. The data is acceptable as no negative impact on data is expected.
DET	Analyte DETECTED
ND	Analyte NOT DETECTED at or above the reporting limit
NR	Not Reported
dry	Sample results reported on a dry weight basis
RPD	Relative Percent Difference

SunStar Laboratories, Inc.

Katherine Shields

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Katherine Shields, Jr. Project Manager

Chain of Custody Record

Date: 4-4-13

Page: 1 OF 1

Project Name: CENCO

Collector: Frane Sosic

Client Project #: 1003-001-300

Batch #: T130794

EDF #:

[illegible]

Sample disposal Instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____

SAMPLE RECEIVING REVIEW SHEET

BATCH # T130794

Client Name: MUREX

Project: CENCO

Received by: PATRICK

Date/Time Received: 4/4/13 16:26

Delivered by: ☐ Client ☒ SunStar Courier ☐ GSO ☐ FedEx ☐ Other _____

Total number of coolers received 0 Temp criteria = 6°C > 0°C (no frozen containers)

Temperature: cooler #1 4.4 °C +/- the CF (- 0.2°C) = 4.2 °C corrected temperature

cooler #2 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

cooler #3 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

Samples outside temp. but received on ice, w/in 6 hours of final sampling. ☒ Yes ☐ No* ☐ N/A

Custody Seals Intact on Cooler/Sample ☐ Yes ☐ No* ☒ N/A

Sample Containers Intact ☒ Yes ☐ No*

Sample labels match COC ID's ☒ Yes ☐ No*

Total number of containers received match COC ☒ Yes ☐ No*

Proper containers received for analyses requested on COC ☒ Yes ☐ No*

Proper preservative indicated on COC/containers for analyses requested ☒ Yes ☐ No* ☐ N/A

Complete shipment received in good condition with correct temperatures, containers, labels, volumes, preservatives and within method specified holding times. ☒ Yes ☐ No*

* Complete Non-Conformance Receiving Sheet if checked

Cooler/Sample Review - Initials and date BC 4/5/13

Comments:



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10 April 2013

Jeremy Squire
Murex
15375 Barranca Parkway, Suite K-101
Irvine, CA 92861
RE: Cenco

Enclosed are the results of analyses for samples received by the laboratory on 04/05/13 16:00. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Wendy Hsiao For Katherine Shields
Jr. Project Manager



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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/10/13 16:41
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ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
LL_W9_040513	T130806-01	Water	04/05/13 10:12	04/05/13 16:00
LL_104A_040513	T130806-02	Water	04/05/13 13:00	04/05/13 16:00
LL_W11_040513	T130806-03	Water	04/05/13 15:00	04/05/13 16:00
LL_TB_040513	T130806-04	Water	04/05/13 00:00	04/05/13 16:00

SunStar Laboratories, Inc.

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Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:41

LL_W9_040513
T130806-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	ND	50	ug/l	1	3040808	04/08/13	04/08/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		119 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040807	04/08/13	04/08/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	1.1	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Wendy Hsiao

Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:41

LL_W9_040513
T130806-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040807	04/08/13	04/08/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

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Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:41

LL_W9_040513
T130806-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3040807	04/08/13	04/08/13	EPA 8260B
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		118 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		101 %	81-136		"	"	"	"
Surrogate: Toluene-d8		95.2 %	88.8-117		"	"	"	"

SunStar Laboratories, Inc.

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/10/13 16:41
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LL_104A_040513
T130806-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	ND	50	ug/l	1	3040808	04/08/13	04/08/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		108 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040807	04/08/13	04/08/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	4.4	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

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Wendy Hsiao

Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:41

LL_104A_040513
T130806-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040807	04/08/13	04/08/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	1.0	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:41

LL_104A_040513
T130806-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3040807	04/08/13	04/08/13	EPA 8260B
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		110 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		98.4 %	81-136		"	"	"	"
Surrogate: Toluene-d8		97.1 %	88.8-117		"	"	"	"

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:41

LL_W11_040513
T130806-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	250	50	ug/l	1	3040808	04/08/13	04/08/13	EPA 8015C
Surrogate: 4-Bromofluorobenzene	115 %	65-135	"	"	"	"	"	"

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040807	04/08/13	04/08/13	EPA 8260B
Bromochloromethane	ND	1.0	"	"	"	"	"	"
Bromodichloromethane	ND	1.0	"	"	"	"	"	"
Bromoform	ND	1.0	"	"	"	"	"	"
Bromomethane	ND	1.0	"	"	"	"	"	"
n-Butylbenzene	ND	1.0	"	"	"	"	"	"
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"
Chlorobenzene	ND	1.0	"	"	"	"	"	"
Chloroethane	ND	1.0	"	"	"	"	"	"
Chloroform	ND	1.0	"	"	"	"	"	"
Chloromethane	ND	1.0	"	"	"	"	"	"
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"
Dibromochloromethane	ND	1.0	"	"	"	"	"	"
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"
Dibromomethane	ND	1.0	"	"	"	"	"	"
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:41

LL_W11_040513
T130806-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040807	04/08/13	04/08/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	8.9	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	1.4	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	14	0.50	"	"	"	"	"	"
Toluene	0.75	0.50	"	"	"	"	"	"
Ethylbenzene	1.2	0.50	"	"	"	"	"	"
m,p-Xylene	3.2	1.0	"	"	"	"	"	"
o-Xylene	0.57	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:41

LL_W11_040513
T130806-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3040807	04/08/13	04/08/13	EPA 8260B
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		119 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		100 %	81-136		"	"	"	"
Surrogate: Toluene-d8		97.4 %	88.8-117		"	"	"	"

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Project: Cenco
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Project Manager: Jeremy Squire

Reported:
04/10/13 16:41

LL_TB_040513
T130806-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040807	04/08/13	04/08/13	EPA 8260B
Bromochloromethane	ND	1.0	"	"	"	"	"	"
Bromodichloromethane	ND	1.0	"	"	"	"	"	"
Bromoform	ND	1.0	"	"	"	"	"	"
Bromomethane	ND	1.0	"	"	"	"	"	"
n-Butylbenzene	ND	1.0	"	"	"	"	"	"
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"
Chlorobenzene	ND	1.0	"	"	"	"	"	"
Chloroethane	ND	1.0	"	"	"	"	"	"
Chloroform	ND	1.0	"	"	"	"	"	"
Chloromethane	ND	1.0	"	"	"	"	"	"
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"
Dibromochloromethane	ND	1.0	"	"	"	"	"	"
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"
Dibromomethane	ND	1.0	"	"	"	"	"	"
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:41

LL_TB_040513
T130806-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

cis-1,3-Dichloropropene	ND	0.50	ug/l	1	3040807	04/08/13	04/08/13	EPA 8260B
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"
Tert-butyl alcohol	ND	10	"	"	"	"	"	"
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"

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Project: Cenco
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Project Manager: Jeremy Squire

Reported:
04/10/13 16:41

LL_TB_040513
T130806-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	ug/l	1	3040807	04/08/13	04/08/13	EPA 8260B	
Surrogate: 4-Bromofluorobenzene	111 %	83.5-119			"	"	"	"	
Surrogate: Dibromofluoromethane	98.5 %	81-136			"	"	"	"	
Surrogate: Toluene-d8	97.9 %	88.8-117			"	"	"	"	

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:41

Purgeable Petroleum Hydrocarbons by EPA 8015C - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040808 - EPA 5030 GC

Blank (3040808-BLK1)

Prepared & Analyzed: 04/08/13

C6-C12 (GRO)	ND	50	ug/l							
Surrogate 4-Bromofluorobenzene	98.3		"	100		98.3	65-135			

LCS (3040808-BS1)

Prepared & Analyzed: 04/08/13

C6-C12 (GRO)	5070	50	ug/l	5520		91.9	75-125			
Surrogate 4-Bromofluorobenzene	82.2		"	100		82.2	65-135			

LCS Dup (3040808-BSD1)

Prepared & Analyzed: 04/08/13

C6-C12 (GRO)	5190	50	ug/l	5520		94.1	75-125	2.36	20	
Surrogate 4-Bromofluorobenzene	76.2		"	100		76.2	65-135			

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:41

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040807 - EPA 5030 GCMS

Blank (3040807-BLK1)

Prepared & Analyzed: 04/08/13

Bromobenzene	ND	1.0	ug/l
Bromochloromethane	ND	1.0	"
Bromodichloromethane	ND	1.0	"
Bromoform	ND	1.0	"
Bromomethane	ND	1.0	"
n-Butylbenzene	ND	1.0	"
sec-Butylbenzene	ND	1.0	"
tert-Butylbenzene	ND	1.0	"
Carbon tetrachloride	ND	0.50	"
Chlorobenzene	ND	1.0	"
Chloroethane	ND	1.0	"
Chloroform	ND	1.0	"
Chloromethane	ND	1.0	"
2-Chlorotoluene	ND	1.0	"
4-Chlorotoluene	ND	1.0	"
Dibromochloromethane	ND	1.0	"
1,2-Dibromo-3-chloropropane	ND	1.0	"
1,2-Dibromoethane (EDB)	ND	1.0	"
Dibromomethane	ND	1.0	"
1,2-Dichlorobenzene	ND	1.0	"
1,3-Dichlorobenzene	ND	1.0	"
1,4-Dichlorobenzene	ND	1.0	"
Dichlorodifluoromethane	ND	0.50	"
1,1-Dichloroethane	ND	1.0	"
1,2-Dichloroethane	ND	0.50	"
1,1-Dichloroethene	ND	1.0	"
cis-1,2-Dichloroethene	ND	1.0	"
trans-1,2-Dichloroethene	ND	1.0	"
1,2-Dichloropropane	ND	1.0	"
1,3-Dichloropropane	ND	1.0	"
2,2-Dichloropropane	ND	1.0	"
1,1-Dichloropropene	ND	1.0	"
cis-1,3-Dichloropropene	ND	0.50	"
trans-1,3-Dichloropropene	ND	0.50	"
Hexachlorobutadiene	ND	1.0	"
Isopropylbenzene	ND	1.0	"

SunStar Laboratories, Inc.

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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:41

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040807 - EPA 5030 GCMS

Blank (3040807-BLK1)

Prepared & Analyzed: 04/08/13

p-Isopropyltoluene	ND	1.0	ug/l							
Methylene chloride	ND	1.0	"							
Naphthalene	ND	1.0	"							
n-Propylbenzene	ND	1.0	"							
Styrene	ND	1.0	"							
1,1,2,2-Tetrachloroethane	ND	1.0	"							
1,1,1,2-Tetrachloroethane	ND	1.0	"							
Tetrachloroethene	ND	1.0	"							
1,2,3-Trichlorobenzene	ND	1.0	"							
1,2,4-Trichlorobenzene	ND	1.0	"							
1,1,2-Trichloroethane	ND	1.0	"							
1,1,1-Trichloroethane	ND	1.0	"							
Trichloroethene	ND	1.0	"							
Trichlorofluoromethane	ND	1.0	"							
1,2,3-Trichloropropane	ND	1.0	"							
1,3,5-Trimethylbenzene	ND	1.0	"							
1,2,4-Trimethylbenzene	ND	1.0	"							
Vinyl chloride	ND	1.0	"							
Benzene	ND	0.50	"							
Toluene	ND	0.50	"							
Ethylbenzene	ND	0.50	"							
m,p-Xylene	ND	1.0	"							
o-Xylene	ND	0.50	"							
Tert-amyl methyl ether	ND	2.0	"							
Tert-butyl alcohol	ND	10	"							
Di-isopropyl ether	ND	2.0	"							
Ethyl tert-butyl ether	ND	2.0	"							
Methyl tert-butyl ether	ND	1.0	"							
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"							
Surrogate 4-Bromofluorobenzene	7.86		"	8.00		98.2	83.5-119			
Surrogate Dibromofluoromethane	6.45		"	8.00		80.6	81-136			S-GC
Surrogate Toluene-d8	7.52		"	8.00		94.0	88.8-117			

SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Wendy Hsiao

Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex	Project: Cenco	Reported:
15375 Barranca Parkway, Suite K-101	Project Number: 1003-001-300	04/10/13 16:41
Irvine CA, 92861	Project Manager: Jeremy Squire	

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040807 - EPA 5030 GCMS

LCS (3040807-BS1)

Prepared: 04/08/13 Analyzed: 04/09/13

Chlorobenzene	21.1	1.0	ug/l	20.0		106	75-125			
1,1-Dichloroethene	21.1	1.0	"	20.0		106	75-125			
Trichloroethene	21.0	1.0	"	20.0		105	75-125			
Benzene	20.7	0.50	"	20.0		103	75-125			
Toluene	21.1	0.50	"	20.0		106	75-125			
Surrogate 4-Bromofluorobenzene	8.67		"	8.00		108	83.5-119			
Surrogate Dibromofluoromethane	7.71		"	8.00		96.4	81-136			
Surrogate Toluene-d8	8.45		"	8.00		106	88.8-117			

LCS Dup (3040807-BSD1)

Prepared & Analyzed: 04/08/13

Chlorobenzene	19.2	1.0	ug/l	20.0		95.8	75-125	9.83	20	
1,1-Dichloroethene	24.8	1.0	"	20.0		124	75-125	16.2	20	
Trichloroethene	21.2	1.0	"	20.0		106	75-125	1.14	20	
Benzene	22.3	0.50	"	20.0		112	75-125	7.77	20	
Toluene	20.3	0.50	"	20.0		101	75-125	4.01	20	
Surrogate 4-Bromofluorobenzene	7.73		"	8.00		96.6	83.5-119			
Surrogate Dibromofluoromethane	8.15		"	8.00		102	81-136			
Surrogate Toluene-d8	7.85		"	8.00		98.1	88.8-117			

SunStar Laboratories, Inc.

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Wendy Hsiao

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/10/13 16:41
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Notes and Definitions

S-GC	Surrogate recovery outside of established control limits. The data was accepted based on valid recovery of the remaining surrogate(s).
DET	Analyte DETECTED
ND	Analyte NOT DETECTED at or above the reporting limit
NR	Not Reported
dry	Sample results reported on a dry weight basis
RPD	Relative Percent Difference

SunStar Laboratories, Inc.

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Wendy Hsiao For Katherine Shields, Jr. Project Manager

Chain of Custody Record

Date: 4-5-13
Project Name: CENCO
Collector: Frane Sosic
Batch #: 7130 806

Page: 1 OF 1

Client Project #: 1003-001-300

EDF #:

Sample ID	Date Sampled	Time	Sample Type	TPHg (8015 M)	VOCs (8260 B)												Total # of containers	Comments/Preservative	Laboratory ID #
LL-W9-040513	4-5-13	10:12	GW	X	X												6		01
LL-104A-040513	4-5-13	1300	GW	X	X												6		02
LL-W11-040513	4-5-13	1800	GW	X	X												6		03
LL-TB-040513			Water		X												2		04
Relinquished by: (signature) 	Date / Time F. Sosic 4-5-2013 16:00	Received by: (Sign / Date / Time) 4/5/13 16:00		Total # of containers				Chain of Custody seals				Notes							
Relinquished by: (signature) 	Date / Time	Received by: (Sign / Date / Time)		Seals intact? Y/N/NA				Received good condition/cold		3.3									
Relinquished by: (signature)	Date / Time	Received by: (Sign / Date / Time)		Turn around time:		Standard													

Sample disposal Instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____

SAMPLE RECEIVING REVIEW SHEET

BATCH # T130806

Client Name: MUREX

Project: CENCO

Received by: PATRICK

Date/Time Received: 4.5.13 / 16:00

Delivered by: ☐ Client ☒ SunStar Courier ☐ GSO ☐ FedEx ☐ Other _____

Total number of coolers received 0 Temp criteria = 6°C > 0°C (no frozen containers)

Temperature: cooler #1 3.5 °C +/- the CF (- 0.2°C) = 3.3 °C corrected temperature

cooler #2 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

cooler #3 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

Samples outside temp. but received on ice, w/in 6 hours of final sampling. ☒ Yes ☐ No* ☐ N/A

Custody Seals Intact on Cooler/Sample ☐ Yes ☐ No* ☒ N/A

Sample Containers Intact ☒ Yes ☐ No*

Sample labels match COC ID's ☒ Yes ☐ No*

Total number of containers received match COC ☒ Yes ☐ No*

Proper containers received for analyses requested on COC ☒ Yes ☐ No*

Proper preservative indicated on COC/containers for analyses requested ☒ Yes ☐ No* ☐ N/A

Complete shipment received in good condition with correct temperatures, containers, labels, volumes preservatives and within method specified holding times. ☒ Yes ☐ No*

* Complete Non-Conformance Receiving Sheet if checked

Cooler/Sample Review - Initials and date SL 4.5.13

Comments:



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10 April 2013

Jeremy Squire
Murex
15375 Barranca Parkway, Suite K-101
Irvine, CA 92861
RE: Cenco

Enclosed are the results of analyses for samples received by the laboratory on 04/08/13 16:00. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Wendy Hsiao For Katherine Shields
Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:05

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
LL_W4_040813	T130812-01	Water	04/08/13 12:00	04/08/13 16:00
LL_503B_040813_01	T130812-02	Water	04/08/13 14:10	04/08/13 16:00
LL_503B_040813_02	T130812-03	Water	04/08/13 14:20	04/08/13 16:00
LL_707_040813	T130812-04	Water	04/08/13 15:43	04/08/13 16:00
LL_TB_040813	T130812-05	Water	04/08/13 00:00	04/08/13 16:00

SunStar Laboratories, Inc.

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Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:05

LL_W4_040813
T130812-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	360	50	ug/l	1	3040919	04/09/13	04/10/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		113 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040918	04/09/13	04/09/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	2.3	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

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15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:05

LL_W4_040813
T130812-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040918	04/09/13	04/09/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	3.8	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	3.8	1.0	"	"	"	"	"	"
Benzene	18	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

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Wendy Hsiao

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15375 Barranca Parkway, Suite K-101
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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:05

LL_W4_040813
T130812-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	77	10	ug/l	1	3040918	04/09/13	04/09/13	EPA 8260B
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	2.8	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene	108 %	83.5-119	"	"	"	"	"	"
Surrogate: Dibromofluoromethane	98.8 %	81-136	"	"	"	"	"	"
Surrogate: Toluene-d8	107 %	88.8-117	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/10/13 16:05
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LL_503B_040813_01
T130812-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	720	50	ug/l	1	3040919	04/09/13	04/10/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene	99.2 %	65-135	"	"	"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040918	04/09/13	04/09/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:05

LL_503B_040813_01
T130812-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040918	04/09/13	04/09/13	EPA 8260B	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	1.4	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	76	1.0	"	"	"	"	"	"	
n-Propylbenzene	4.6	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	8.7	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	39	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	64	0.50	"	"	"	"	"	"	
Toluene	4.3	0.50	"	"	"	"	"	"	
Ethylbenzene	17	0.50	"	"	"	"	"	"	
m,p-Xylene	47	1.0	"	"	"	"	"	"	
o-Xylene	12	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	20	10	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:05

LL_503B_040813_01
T130812-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Di-isopropyl ether	ND	2.0	ug/l	1	3040918	04/09/13	04/09/13	EPA 8260B
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	2.8	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		111 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		99.9 %	81-136		"	"	"	"
Surrogate: Toluene-d8		105 %	88.8-117		"	"	"	"

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:05

LL_503B_040813_02
T130812-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	600	50	ug/l	1	3040919	04/09/13	04/10/13	EPA 8015C
Surrogate: 4-Bromofluorobenzene	94.0 %	65-135	"	"	"	"	"	"

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040918	04/09/13	04/09/13	EPA 8260B
Bromochloromethane	ND	1.0	"	"	"	"	"	"
Bromodichloromethane	ND	1.0	"	"	"	"	"	"
Bromoform	ND	1.0	"	"	"	"	"	"
Bromomethane	ND	1.0	"	"	"	"	"	"
n-Butylbenzene	ND	1.0	"	"	"	"	"	"
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"
Chlorobenzene	ND	1.0	"	"	"	"	"	"
Chloroethane	ND	1.0	"	"	"	"	"	"
Chloroform	ND	1.0	"	"	"	"	"	"
Chloromethane	ND	1.0	"	"	"	"	"	"
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"
Dibromochloromethane	ND	1.0	"	"	"	"	"	"
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"
Dibromomethane	ND	1.0	"	"	"	"	"	"
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:05

LL_503B_040813_02
T130812-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040918	04/09/13	04/09/13	EPA 8260B	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	1.3	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	75	1.0	"	"	"	"	"	"	
n-Propylbenzene	4.2	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	8.1	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	36	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	62	0.50	"	"	"	"	"	"	
Toluene	4.1	0.50	"	"	"	"	"	"	
Ethylbenzene	16	0.50	"	"	"	"	"	"	
m,p-Xylene	44	1.0	"	"	"	"	"	"	
o-Xylene	11	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	19	10	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:05

LL_503B_040813_02
T130812-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Di-isopropyl ether	ND	2.0	ug/l	1	3040918	04/09/13	04/09/13	EPA 8260B
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	2.7	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		108 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		98.4 %	81-136		"	"	"	"
Surrogate: Toluene-d8		104 %	88.8-117		"	"	"	"

SunStar Laboratories, Inc.

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/10/13 16:05
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LL_707_040813
T130812-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	240	50	ug/l	1	3040919	04/09/13	04/10/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene	97.4 %	65-135	"	"	"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040918	04/09/13	04/09/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:05

LL_707_040813
T130812-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3040918	04/09/13	04/09/13	EPA 8260B	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	1.2	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	29	1.0	"	"	"	"	"	"	
n-Propylbenzene	2.8	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	2.7	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	6.0	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	92	0.50	"	"	"	"	"	"	
Toluene	5.6	0.50	"	"	"	"	"	"	
Ethylbenzene	5.2	0.50	"	"	"	"	"	"	
m,p-Xylene	27	1.0	"	"	"	"	"	"	
o-Xylene	5.0	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	ND	10	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:05

LL_707_040813
T130812-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Di-isopropyl ether	ND	2.0	ug/l	1	3040918	04/09/13	04/09/13	EPA 8260B
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	2.1	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene	110 %	83.5-119	"	"	"	"	"	"
Surrogate: Dibromofluoromethane	99.2 %	81-136	"	"	"	"	"	"
Surrogate: Toluene-d8	105 %	88.8-117	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:05

LL_TB_040813
T130812-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3040918	04/09/13	04/09/13	EPA 8260B
Bromochloromethane	ND	1.0	"	"	"	"	"	"
Bromodichloromethane	ND	1.0	"	"	"	"	"	"
Bromoform	ND	1.0	"	"	"	"	"	"
Bromomethane	ND	1.0	"	"	"	"	"	"
n-Butylbenzene	ND	1.0	"	"	"	"	"	"
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"
Chlorobenzene	ND	1.0	"	"	"	"	"	"
Chloroethane	ND	1.0	"	"	"	"	"	"
Chloroform	ND	1.0	"	"	"	"	"	"
Chloromethane	ND	1.0	"	"	"	"	"	"
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"
Dibromochloromethane	ND	1.0	"	"	"	"	"	"
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"
Dibromomethane	ND	1.0	"	"	"	"	"	"
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:05

LL_TB_040813
T130812-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

cis-1,3-Dichloropropene	ND	0.50	ug/l	1	3040918	04/09/13	04/09/13	EPA 8260B	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	ND	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	ND	1.0	"	"	"	"	"	"	
n-Propylbenzene	ND	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	ND	0.50	"	"	"	"	"	"	
Toluene	ND	0.50	"	"	"	"	"	"	
Ethylbenzene	ND	0.50	"	"	"	"	"	"	
m,p-Xylene	ND	1.0	"	"	"	"	"	"	
o-Xylene	ND	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	ND	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Wendy Hsiao

Wendy Hsiao For Katherine Shields, Jr. Project Manager



25712 Commercentre Drive
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949.297.5020 Phone
949.297.5027 Fax

Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:05

LL_TB_040813
T130812-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	ug/l	1	3040918	04/09/13	04/09/13	EPA 8260B	
Surrogate: 4-Bromofluorobenzene	109 %	83.5-119			"	"	"	"	
Surrogate: Dibromofluoromethane	100 %	81-136			"	"	"	"	
Surrogate: Toluene-d8	103 %	88.8-117			"	"	"	"	

SunStar Laboratories, Inc.

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Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/10/13 16:05
--	--	-----------------------------

Purgeable Petroleum Hydrocarbons by EPA 8015C - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
Batch 3040919 - EPA 5030 GC										
Blank (3040919-BLK1)				Prepared: 04/09/13 Analyzed: 04/10/13						
C6-C12 (GRO)	ND	50	ug/l							
Surrogate 4-Bromofluorobenzene	107		"	100		107	65-135			
LCS (3040919-BS1)				Prepared: 04/09/13 Analyzed: 04/10/13						
C6-C12 (GRO)	5360	50	ug/l	5520		97.0	75-125			
Surrogate 4-Bromofluorobenzene	73.6		"	100		73.6	65-135			
Matrix Spike (3040919-MS1)				Source: T130812-04		Prepared: 04/09/13 Analyzed: 04/10/13				
C6-C12 (GRO)	5010	50	ug/l	5520	242	86.4	65-135			
Surrogate 4-Bromofluorobenzene	82.7		"	100		82.7	65-135			
Matrix Spike Dup (3040919-MSD1)				Source: T130812-04		Prepared: 04/09/13 Analyzed: 04/10/13				
C6-C12 (GRO)	5460	50	ug/l	5520	242	94.5	65-135	8.54	20	
Surrogate 4-Bromofluorobenzene	89.8		"	100		89.8	65-135			

SunStar Laboratories, Inc.

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Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:05

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040918 - EPA 5030 GCMS

Blank (3040918-BLK1)

Prepared & Analyzed: 04/09/13

Bromobenzene	ND	1.0	ug/l
Bromochloromethane	ND	1.0	"
Bromodichloromethane	ND	1.0	"
Bromoform	ND	1.0	"
Bromomethane	ND	1.0	"
n-Butylbenzene	ND	1.0	"
sec-Butylbenzene	ND	1.0	"
tert-Butylbenzene	ND	1.0	"
Carbon tetrachloride	ND	0.50	"
Chlorobenzene	ND	1.0	"
Chloroethane	ND	1.0	"
Chloroform	ND	1.0	"
Chloromethane	ND	1.0	"
2-Chlorotoluene	ND	1.0	"
4-Chlorotoluene	ND	1.0	"
Dibromochloromethane	ND	1.0	"
1,2-Dibromo-3-chloropropane	ND	1.0	"
1,2-Dibromoethane (EDB)	ND	1.0	"
Dibromomethane	ND	1.0	"
1,2-Dichlorobenzene	ND	1.0	"
1,3-Dichlorobenzene	ND	1.0	"
1,4-Dichlorobenzene	ND	1.0	"
Dichlorodifluoromethane	ND	0.50	"
1,1-Dichloroethane	ND	1.0	"
1,2-Dichloroethane	ND	0.50	"
1,1-Dichloroethene	ND	1.0	"
cis-1,2-Dichloroethene	ND	1.0	"
trans-1,2-Dichloroethene	ND	1.0	"
1,2-Dichloropropane	ND	1.0	"
1,3-Dichloropropane	ND	1.0	"
2,2-Dichloropropane	ND	1.0	"
1,1-Dichloropropene	ND	1.0	"
cis-1,3-Dichloropropene	ND	0.50	"
trans-1,3-Dichloropropene	ND	0.50	"
Hexachlorobutadiene	ND	1.0	"
Isopropylbenzene	ND	1.0	"

SunStar Laboratories, Inc.

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Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/10/13 16:05

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040918 - EPA 5030 GCMS

Blank (3040918-BLK1)

Prepared & Analyzed: 04/09/13

p-Isopropyltoluene	ND	1.0	ug/l
Methylene chloride	ND	1.0	"
Naphthalene	ND	1.0	"
n-Propylbenzene	ND	1.0	"
Styrene	ND	1.0	"
1,1,2,2-Tetrachloroethane	ND	1.0	"
1,1,1,2-Tetrachloroethane	ND	1.0	"
Tetrachloroethene	ND	1.0	"
1,2,3-Trichlorobenzene	ND	1.0	"
1,2,4-Trichlorobenzene	ND	1.0	"
1,1,2-Trichloroethane	ND	1.0	"
1,1,1-Trichloroethane	ND	1.0	"
Trichloroethene	ND	1.0	"
Trichlorofluoromethane	ND	1.0	"
1,2,3-Trichloropropane	ND	1.0	"
1,3,5-Trimethylbenzene	ND	1.0	"
1,2,4-Trimethylbenzene	ND	1.0	"
Vinyl chloride	ND	1.0	"
Benzene	ND	0.50	"
Toluene	ND	0.50	"
Ethylbenzene	ND	0.50	"
m,p-Xylene	ND	1.0	"
o-Xylene	ND	0.50	"
Tert-amyl methyl ether	ND	2.0	"
Tert-butyl alcohol	ND	10	"
Di-isopropyl ether	ND	2.0	"
Ethyl tert-butyl ether	ND	2.0	"
Methyl tert-butyl ether	ND	1.0	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"
Surrogate 4-Bromofluorobenzene	8.79		"
Surrogate Dibromofluoromethane	8.02		"
Surrogate Toluene-d8	8.27		"

8.00 110 83.5-119
8.00 100 81-136
8.00 103 88.8-117

SunStar Laboratories, Inc.

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Wendy Hsiao

Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex	Project: Cenco	Reported:
15375 Barranca Parkway, Suite K-101	Project Number: 1003-001-300	04/10/13 16:05
Irvine CA, 92861	Project Manager: Jeremy Squire	

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3040918 - EPA 5030 GCMS

LCS (3040918-BS1)

Prepared & Analyzed: 04/09/13

Chlorobenzene	17.6	1.0	ug/l	20.0		88.2	75-125			
1,1-Dichloroethene	19.7	1.0	"	20.0		98.6	75-125			
Trichloroethene	17.7	1.0	"	20.0		88.6	75-125			
Benzene	19.3	0.50	"	20.0		96.4	75-125			
Toluene	15.8	0.50	"	20.0		79.2	75-125			
Surrogate 4-Bromofluorobenzene	8.30		"	8.00		104	83.5-119			
Surrogate Dibromofluoromethane	8.60		"	8.00		108	81-136			
Surrogate Toluene-d8	8.10		"	8.00		101	88.8-117			

Matrix Spike (3040918-MS1)

Source: T130812-01

Prepared & Analyzed: 04/09/13

Chlorobenzene	16.5	1.0	ug/l	20.0	ND	82.5	75-125			
1,1-Dichloroethene	19.8	1.0	"	20.0	ND	99.2	75-125			
Trichloroethene	18.2	1.0	"	20.0	ND	90.8	75-125			
Benzene	38.5	0.50	"	20.0	18.1	102	75-125			
Toluene	16.3	0.50	"	20.0	ND	81.4	75-125			
Surrogate 4-Bromofluorobenzene	8.10		"	8.00		101	83.5-119			
Surrogate Dibromofluoromethane	8.45		"	8.00		106	81-136			
Surrogate Toluene-d8	8.25		"	8.00		103	88.8-117			

Matrix Spike Dup (3040918-MSD1)

Source: T130812-01

Prepared & Analyzed: 04/09/13

Chlorobenzene	17.1	1.0	ug/l	20.0	ND	85.4	75-125	3.40	20	
1,1-Dichloroethene	19.2	1.0	"	20.0	ND	96.1	75-125	3.22	20	
Trichloroethene	17.8	1.0	"	20.0	ND	88.8	75-125	2.23	20	
Benzene	38.7	0.50	"	20.0	18.1	103	75-125	0.544	20	
Toluene	16.3	0.50	"	20.0	ND	81.4	75-125	0.123	20	
Surrogate 4-Bromofluorobenzene	8.12		"	8.00		102	83.5-119			
Surrogate Dibromofluoromethane	8.38		"	8.00		105	81-136			
Surrogate Toluene-d8	8.12		"	8.00		102	88.8-117			

SunStar Laboratories, Inc.

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Wendy Hsiao

Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/10/13 16:05
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Notes and Definitions

DET Analyte DETECTED
ND Analyte NOT DETECTED at or above the reporting limit
NR Not Reported
dry Sample results reported on a dry weight basis
RPD Relative Percent Difference

SunStar Laboratories, Inc.

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Wendy Hsiao For Katherine Shields, Jr. Project Manager

Chain of Custody Record

Date: 4-8-13
Project Name: CENCO
Collector: Frane Sosic
Batch #: T130812

Page: 1 OF 1

Client Project #: 1003-001-300

EDF #:

[illegible]

Sample disposal Instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____

SAMPLE RECEIVING REVIEW SHEET

BATCH # T130812

Client Name: Murex

Project: Cenco

Received by: Patrick

Date/Time Received: 4/8/13 1600

Delivered by: ☐ Client ☒ SunStar Courier ☐ GSO ☐ FedEx ☐ Other _____

Total number of coolers received 1 Temp criteria = 6°C > 0°C (no frozen containers)

Temperature: cooler #1 4.0 °C +/- the CF (- 0.2°C) = 3.8 °C corrected temperature

cooler #2 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

cooler #3 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

Samples outside temp. but received on ice, w/in 6 hours of final sampling. ☒ Yes ☐ No* ☐ N/A

Custody Seals Intact on Cooler/Sample ☐ Yes ☐ No* ☒ N/A

Sample Containers Intact ☒ Yes ☐ No*

Sample labels match COC ID's ☒ Yes ☐ No*

Total number of containers received match COC ☒ Yes ☐ No*


Proper containers received for analyses requested on COC ☒ Yes ☐ No*

Proper preservative indicated on COC/containers for analyses requested ☒ Yes ☐ No* ☐ N/A

Complete shipment received in good condition with correct temperatures, containers, labels, volumes preservatives and within method specified holding times. ☒ Yes ☐ No*

* Complete Non-Conformance Receiving Sheet if checked

Cooler/Sample Review - Initials and date

 4/8/13

Comments:



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12 April 2013

Jeremy Squire
Murex
15375 Barranca Parkway, Suite K-101
Irvine, CA 92861
RE: Cenco

Enclosed are the results of analyses for samples received by the laboratory on 04/09/13 16:36. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Katherine Shields
Jr. Project Manager



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949.297.5027 Fax

Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/12/13 13:52

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
LL_17A_040913	T130820-01	Water	04/09/13 11:11	04/09/13 16:36
LL_17B_040913	T130820-02	Water	04/09/13 13:10	04/09/13 16:36
LL_17C_040913	T130820-03	Water	04/09/13 16:12	04/09/13 16:36
LL_TB_040913	T130820-04	Water	04/09/13 00:00	04/09/13 16:36

SunStar Laboratories, Inc.

Katherine Shields

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Katherine Shields, Jr. Project Manager



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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/12/13 13:52

LL_17A_040913
T130820-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	ND	50	ug/l	1	3041011	04/10/13	04/11/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		97.1 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041012	04/10/13	04/10/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager

Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/12/13 13:52

LL_17A_040913
T130820-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3041012	04/10/13	04/10/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/12/13 13:52

LL_17A_040913
T130820-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3041012	04/10/13	04/10/13	EPA 8260B
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		106 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		90.2 %	81-136		"	"	"	"
Surrogate: Toluene-d8		99.4 %	88.8-117		"	"	"	"

SunStar Laboratories, Inc.

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/12/13 13:52
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LL_17B_040913
T130820-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	ND	50	ug/l	1	3041011	04/10/13	04/11/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		100 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041012	04/10/13	04/10/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/12/13 13:52

LL_17B_040913
T130820-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3041012	04/10/13	04/10/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/12/13 13:52

LL_17B_040913
T130820-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3041012	04/10/13	04/10/13	EPA 8260B
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene	106 %	83.5-119	"	"	"	"	"	"
Surrogate: Dibromofluoromethane	90.6 %	81-136	"	"	"	"	"	"
Surrogate: Toluene-d8	96.0 %	88.8-117	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/12/13 13:52

LL_17C_040913
T130820-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	ND	50	ug/l	1	3041011	04/10/13	04/11/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		108 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041012	04/10/13	04/10/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/12/13 13:52

LL_17C_040913
T130820-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3041012	04/10/13	04/10/13	EPA 8260B	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	ND	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	ND	1.0	"	"	"	"	"	"	
n-Propylbenzene	ND	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	ND	0.50	"	"	"	"	"	"	
Toluene	ND	0.50	"	"	"	"	"	"	
Ethylbenzene	ND	0.50	"	"	"	"	"	"	
m,p-Xylene	ND	1.0	"	"	"	"	"	"	
o-Xylene	ND	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/12/13 13:52

LL_17C_040913
T130820-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3041012	04/10/13	04/10/13	EPA 8260B
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		103 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		94.6 %	81-136		"	"	"	"
Surrogate: Toluene-d8		98.6 %	88.8-117		"	"	"	"

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/12/13 13:52

LL_TB_040913
T130820-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041012	04/10/13	04/10/13	EPA 8260B
Bromochloromethane	ND	1.0	"	"	"	"	"	"
Bromodichloromethane	ND	1.0	"	"	"	"	"	"
Bromoform	ND	1.0	"	"	"	"	"	"
Bromomethane	ND	1.0	"	"	"	"	"	"
n-Butylbenzene	ND	1.0	"	"	"	"	"	"
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"
Chlorobenzene	ND	1.0	"	"	"	"	"	"
Chloroethane	ND	1.0	"	"	"	"	"	"
Chloroform	ND	1.0	"	"	"	"	"	"
Chloromethane	ND	1.0	"	"	"	"	"	"
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"
Dibromochloromethane	ND	1.0	"	"	"	"	"	"
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"
Dibromomethane	ND	1.0	"	"	"	"	"	"
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/12/13 13:52

LL_TB_040913
T130820-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

cis-1,3-Dichloropropene	ND	0.50	ug/l	1	3041012	04/10/13	04/10/13	EPA 8260B
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"
Tert-butyl alcohol	ND	10	"	"	"	"	"	"
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Lake Forest, California 92630
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949.297.5027 Fax

Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/12/13 13:52

LL_TB_040913
T130820-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	ug/l	1	3041012	04/10/13	04/10/13	EPA 8260B	
Surrogate: 4-Bromofluorobenzene	104 %	83.5-119			"	"	"	"	
Surrogate: Dibromofluoromethane	96.2 %	81-136			"	"	"	"	
Surrogate: Toluene-d8	94.5 %	88.8-117			"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/12/13 13:52
--	--	-----------------------------

Purgeable Petroleum Hydrocarbons by EPA 8015C - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
Batch 3041011 - EPA 5030 GC										
Blank (3041011-BLK1)				Prepared: 04/10/13 Analyzed: 04/11/13						
C6-C12 (GRO)	ND	50	ug/l							
Surrogate 4-Bromofluorobenzene	89.8		"	100		89.8	65-135			
LCS (3041011-BS1)				Prepared: 04/10/13 Analyzed: 04/11/13						
C6-C12 (GRO)	5110	50	ug/l	5520		92.6	75-125			
Surrogate 4-Bromofluorobenzene	80.2		"	100		80.2	65-135			
Matrix Spike (3041011-MS1)				Source: T130820-01		Prepared: 04/10/13 Analyzed: 04/11/13				
C6-C12 (GRO)	5000	50	ug/l	5520	27.7	90.1	65-135			
Surrogate 4-Bromofluorobenzene	75.4		"	100		75.4	65-135			
Matrix Spike Dup (3041011-MSD1)				Source: T130820-01		Prepared: 04/10/13 Analyzed: 04/11/13				
C6-C12 (GRO)	4960	50	ug/l	5520	27.7	89.4	65-135	0.853	20	
Surrogate 4-Bromofluorobenzene	83.1		"	100		83.1	65-135			

SunStar Laboratories, Inc.

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Katherine Shields

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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/12/13 13:52

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041012 - EPA 5030 GCMS

Blank (3041012-BLK1)

Prepared & Analyzed: 04/10/13

Bromobenzene	ND	1.0	ug/l
Bromochloromethane	ND	1.0	"
Bromodichloromethane	ND	1.0	"
Bromoform	ND	1.0	"
Bromomethane	ND	1.0	"
n-Butylbenzene	ND	1.0	"
sec-Butylbenzene	ND	1.0	"
tert-Butylbenzene	ND	1.0	"
Carbon tetrachloride	ND	0.50	"
Chlorobenzene	ND	1.0	"
Chloroethane	ND	1.0	"
Chloroform	ND	1.0	"
Chloromethane	ND	1.0	"
2-Chlorotoluene	ND	1.0	"
4-Chlorotoluene	ND	1.0	"
Dibromochloromethane	ND	1.0	"
1,2-Dibromo-3-chloropropane	ND	1.0	"
1,2-Dibromoethane (EDB)	ND	1.0	"
Dibromomethane	ND	1.0	"
1,2-Dichlorobenzene	ND	1.0	"
1,3-Dichlorobenzene	ND	1.0	"
1,4-Dichlorobenzene	ND	1.0	"
Dichlorodifluoromethane	ND	0.50	"
1,1-Dichloroethane	ND	1.0	"
1,2-Dichloroethane	ND	0.50	"
1,1-Dichloroethene	ND	1.0	"
cis-1,2-Dichloroethene	ND	1.0	"
trans-1,2-Dichloroethene	ND	1.0	"
1,2-Dichloropropane	ND	1.0	"
1,3-Dichloropropane	ND	1.0	"
2,2-Dichloropropane	ND	1.0	"
1,1-Dichloropropene	ND	1.0	"
cis-1,3-Dichloropropene	ND	0.50	"
trans-1,3-Dichloropropene	ND	0.50	"
Hexachlorobutadiene	ND	1.0	"
Isopropylbenzene	ND	1.0	"

SunStar Laboratories, Inc.

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Katherine Shields

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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/12/13 13:52

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041012 - EPA 5030 GCMS

Blank (3041012-BLK1)

Prepared & Analyzed: 04/10/13

p-Isopropyltoluene	ND	1.0	ug/l							
Methylene chloride	ND	1.0	"							
Naphthalene	ND	1.0	"							
n-Propylbenzene	ND	1.0	"							
Styrene	ND	1.0	"							
1,1,2,2-Tetrachloroethane	ND	1.0	"							
1,1,1,2-Tetrachloroethane	ND	1.0	"							
Tetrachloroethene	ND	1.0	"							
1,2,3-Trichlorobenzene	ND	1.0	"							
1,2,4-Trichlorobenzene	ND	1.0	"							
1,1,2-Trichloroethane	ND	1.0	"							
1,1,1-Trichloroethane	ND	1.0	"							
Trichloroethene	ND	1.0	"							
Trichlorofluoromethane	ND	1.0	"							
1,2,3-Trichloropropane	ND	1.0	"							
1,3,5-Trimethylbenzene	ND	1.0	"							
1,2,4-Trimethylbenzene	ND	1.0	"							
Vinyl chloride	ND	1.0	"							
Benzene	ND	0.50	"							
Toluene	ND	0.50	"							
Ethylbenzene	ND	0.50	"							
m,p-Xylene	ND	1.0	"							
o-Xylene	ND	0.50	"							
Tert-amyl methyl ether	ND	2.0	"							
Tert-butyl alcohol	ND	10	"							
Di-isopropyl ether	ND	2.0	"							
Ethyl tert-butyl ether	ND	2.0	"							
Methyl tert-butyl ether	ND	1.0	"							
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"							
Surrogate 4-Bromofluorobenzene	7.70		"	8.00		96.2	83.5-119			
Surrogate Dibromofluoromethane	6.06		"	8.00		75.8	81-136			S-GC
Surrogate Toluene-d8	7.73		"	8.00		96.6	88.8-117			

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/12/13 13:52

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041012 - EPA 5030 GCMS

LCS (3041012-BS1)

Prepared & Analyzed: 04/10/13

Chlorobenzene	19.1	1.0	ug/l	20.0		95.7	75-125			
1,1-Dichloroethene	24.8	1.0	"	20.0		124	75-125			
Trichloroethene	21.8	1.0	"	20.0		109	75-125			
Benzene	21.2	0.50	"	20.0		106	75-125			
Toluene	19.6	0.50	"	20.0		98.0	75-125			
Surrogate 4-Bromofluorobenzene	8.20		"	8.00		102	83.5-119			
Surrogate Dibromofluoromethane	7.60		"	8.00		95.0	81-136			
Surrogate Toluene-d8	7.61		"	8.00		95.1	88.8-117			

Matrix Spike (3041012-MS1)

Source: T130819-01

Prepared & Analyzed: 04/10/13

Chlorobenzene	19.7	1.0	ug/l	20.0	ND	98.6	75-125			
1,1-Dichloroethene	26.8	1.0	"	20.0	ND	134	75-125			QM-05
Trichloroethene	22.1	1.0	"	20.0	ND	111	75-125			
Benzene	21.4	0.50	"	20.0	ND	107	75-125			
Toluene	20.2	0.50	"	20.0	ND	101	75-125			
Surrogate 4-Bromofluorobenzene	7.59		"	8.00		94.9	83.5-119			
Surrogate Dibromofluoromethane	7.69		"	8.00		96.1	81-136			
Surrogate Toluene-d8	7.87		"	8.00		98.4	88.8-117			

Matrix Spike Dup (3041012-MSD1)

Source: T130819-01

Prepared & Analyzed: 04/10/13

Chlorobenzene	23.0	1.0	ug/l	20.0	ND	115	75-125	15.3	20	
1,1-Dichloroethene	30.0	1.0	"	20.0	ND	150	75-125	11.2	20	QM-05
Trichloroethene	25.7	1.0	"	20.0	ND	128	75-125	14.8	20	QM-05
Benzene	25.3	0.50	"	20.0	ND	126	75-125	16.5	20	QM-05
Toluene	23.7	0.50	"	20.0	ND	118	75-125	15.8	20	
Surrogate 4-Bromofluorobenzene	7.71		"	8.00		96.4	83.5-119			
Surrogate Dibromofluoromethane	7.58		"	8.00		94.8	81-136			
Surrogate Toluene-d8	7.78		"	8.00		97.2	88.8-117			

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/12/13 13:52
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Notes and Definitions

S-GC Surrogate recovery outside of established control limits. The data was accepted based on valid recovery of the remaining surrogate(s).

QM-05 The spike recovery was outside acceptance limits for the MS and/or MSD due to possible matrix interference. The LCS was within acceptance criteria. The data is acceptable as no negative impact on data is expected.

DET Analyte DETECTED

ND Analyte NOT DETECTED at or above the reporting limit

NR Not Reported

dry Sample results reported on a dry weight basis

RPD Relative Percent Difference

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager

Chain of Custody Record

Date: 4-9-13

Page: 1 OF 1

Project Name: CENCO

Collector: Frane Sosic

Client Project #: 1003-001-300

Batch #: T130820

EDF #:

[illegible]

Sample disposal instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____

SAMPLE RECEIVING REVIEW SHEET

BATCH # T130820

Client Name: MUREX

Project: CEND

Received by: PATRICK

Date/Time Received: 4.9.13 / 16:36

Delivered by : ☐ Client ☒ SunStar Courier ☐ GSO ☐ FedEx ☐ Other _____

Total number of coolers received 0 Temp criteria = 6°C > 0°C (no frozen containers)

Temperature: cooler #1 3.3 °C +/- the CF (- 0.2°C) = 3.1 °C corrected temperature

cooler #2 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

cooler #3 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

Samples outside temp. but received on ice, w/in 6 hours of final sampling. ☒ Yes ☐ No* ☐ N/A

Custody Seals Intact on Cooler/Sample ☐ Yes ☐ No* ☒ N/A

Sample Containers Intact ☒ Yes ☐ No*

Sample labels match COC ID's ☒ Yes ☐ No*

Total number of containers received match COC ☒ Yes ☐ No*

Proper containers received for analyses requested on COC ☒ Yes ☐ No*

Proper preservative indicated on COC/containers for analyses requested ☒ Yes ☐ No* ☐ N/A

Complete shipment received in good condition with correct temperatures, containers, labels, volumes preservatives and within method specified holding times. ☒ Yes ☐ No*

* Complete Non-Conformance Receiving Sheet if checked

Cooler/Sample Review - Initials and date SL 4.9.13

Comments:



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17 April 2013

Jeremy Squire
Murex
15375 Barranca Parkway, Suite K-101
Irvine, CA 92861
RE: Cenco

Enclosed are the results of analyses for samples received by the laboratory on 04/10/13 17:00. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Katherine Shields
Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
LL_EW1_041013	T130836-01	Water	04/10/13 10:00	04/10/13 17:00
LL_703_041013	T130836-02	Water	04/10/13 12:45	04/10/13 17:00
LL_701_041013	T130836-03	Water	04/10/13 14:13	04/10/13 17:00
LL_702_041013	T130836-04	Water	04/10/13 15:42	04/10/13 17:00
LL_705_041013	T130836-05	Water	04/10/13 16:36	04/10/13 17:00
LL_TB_041013	T130836-06	Water	04/10/13 00:00	04/10/13 17:00

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

LL_EW1_041013

T130836-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	1400	50	ug/l	1	3041128	04/11/13	04/12/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		182 %	65-135		"	"	"	"	S-04

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041124	04/11/13	04/11/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	5.9	1.0	"	"	"	"	"	"	
sec-Butylbenzene	10	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

LL_EW1_041013
T130836-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3041124	04/11/13	04/11/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	13	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	88	1.0	"	"	"	"	"	"
n-Propylbenzene	17	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	2,1	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

LL_EW1_041013
T130836-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3041124	04/11/13	04/11/13	EPA 8260B
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene	97.6 %	83.5-119	"	"	"	"	"	"
Surrogate: Dibromofluoromethane	110 %	81-136	"	"	"	"	"	"
Surrogate: Toluene-d8	103 %	88.8-117	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

LL_703_041013
T130836-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	ND	50	ug/l	1	3041128	04/11/13	04/12/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		88.4 %	65-135		"	"	"	"	

Metals by SM 3500 Series Methods

Ferrous Iron	ND	0.100	mg/l	1	3041618	04/11/13	04/16/13	EPA6010/S M3500	
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Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041124	04/11/13	04/12/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

LL_703_041013
T130836-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1-Dichloroethene	ND	1.0	ug/l	1	3041124	04/11/13	04/12/13	EPA 8260B
cis-1,2-Dichloroethene	5.7	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

LL_703_041013
T130836-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

m,p-Xylene	ND	1.0	ug/l	1	3041124	04/11/13	04/12/13	EPA 8260B
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"
Tert-butyl alcohol	ND	10	"	"	"	"	"	"
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene	114 %	83.5-119	"	"	"	"	"	"
Surrogate: Dibromofluoromethane	102 %	81-136	"	"	"	"	"	"
Surrogate: Toluene-d8	102 %	88.8-117	"	"	"	"	"	"

Conventional Chemistry Parameters by APHA/EPA/ASTM Methods

Total Alkalinity	560	20	mg/l	1	3041122	04/11/13	04/12/13	EPA 310.1
Anions by EPA Method 300.0								
Sulfate as SO ₄	383	5.00	mg/l	10	3041121	04/11/13	04/11/13	EPA 300.0
Nitrate as NO ₃	ND	0.500	"	1	"	"	04/11/13	"

RSK-175

Methane	1.17	1.00	ug/l	1	3041129	04/11/13	04/12/13	RSK-175
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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/17/13 14:10
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LL_701_041013
T130836-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	ND	50	ug/l	1	3041128	04/11/13	04/12/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		107 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041124	04/11/13	04/12/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	7.3	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

LL_701_041013
T130836-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3041124	04/11/13	04/12/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	1.8	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

LL_701_041013
T130836-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3041124	04/11/13	04/12/13	EPA 8260B
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		110 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		103 %	81-136		"	"	"	"
Surrogate: Toluene-d8		102 %	88.8-117		"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

LL_702_041013
T130836-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	2300	50	ug/l	1	3041128	04/11/13	04/12/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		129 %	65-135		"	"	"	"	

Metals by SM 3500 Series Methods

Ferrous Iron	ND	0.100	mg/l	1	3041618	04/11/13	04/16/13	EPA6010/S M3500	
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Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041124	04/11/13	04/12/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	17	1.0	"	"	"	"	"	"	
tert-Butylbenzene	3.0	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	2.1	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	4.4	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	1.5	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

LL_702_041013
T130836-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1-Dichloroethene	ND	1.0	ug/l	1	3041124	04/11/13	04/12/13	EPA 8260B
cis-1,2-Dichloroethene	1.3	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
1,2-Dichloropropane	13	1.0	"	"	"	"	"	"
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	6.2	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	1.8	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	15	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

LL_702_041013
T130836-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

m,p-Xylene	ND	1.0	ug/l	1	3041124	04/11/13	04/12/13	EPA 8260B
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"
Tert-butyl alcohol	ND	10	"	"	"	"	"	"
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene	110 %	83.5-119	"	"	"	"	"	"
Surrogate: Dibromofluoromethane	104 %	81-136	"	"	"	"	"	"
Surrogate: Toluene-d8	102 %	88.8-117	"	"	"	"	"	"

Conventional Chemistry Parameters by APHA/EPA/ASTM Methods

Total Alkalinity	840	20	mg/l	1	3041122	04/11/13	04/12/13	EPA 310.1
Anions by EPA Method 300.0								
Sulfate as SO ₄	101	5.00	mg/l	10	3041121	04/11/13	04/11/13	EPA 300.0
Nitrate as NO ₃	ND	0.500	"	1	"	"	04/11/13	"

RSK-175

Methane	757	3.00	ug/l	3	3041129	04/11/13	04/12/13	RSK-175
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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/17/13 14:10
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LL_705_041013
T130836-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	140	50	ug/l	1	3041128	04/11/13	04/12/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		104 %	65-135		"	"	"	"	

Metals by SM 3500 Series Methods

Ferrous Iron	ND	0.100	mg/l	1	3041618	04/11/13	04/16/13	EPA6010/S M3500	
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Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041124	04/11/13	04/12/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	1.8	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

LL_705_041013
T130836-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1-Dichloroethene	1.5	1.0	ug/l	1	3041124	04/11/13	04/12/13	EPA 8260B
cis-1,2-Dichloroethene	7.8	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	1.9	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	2.0	1.0	"	"	"	"	"	"
Benzene	0.97	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

LL_705_041013
T130836-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

m,p-Xylene	ND	1.0	ug/l	1	3041124	04/11/13	04/12/13	EPA 8260B	
o-Xylene	ND	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	ND	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	10	1.0	"	"	"	"	"	"	
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene		115 %	83.5-119		"	"	"	"	
Surrogate: Dibromofluoromethane		99.6 %	81-136		"	"	"	"	
Surrogate: Toluene-d8		103 %	88.8-117		"	"	"	"	

Conventional Chemistry Parameters by APHA/EPA/ASTM Methods

Total Alkalinity	640	20	mg/l	1	3041122	04/11/13	04/12/13	EPA 310.1	
Anions by EPA Method 300.0									
Sulfate as SO4	95.6	5.00	mg/l	10	3041121	04/11/13	04/11/13	EPA 300.0	
Nitrate as NO3	ND	0.500	"	1	"	"	04/11/13	"	

RSK-175

Methane	99.3	1.00	ug/l	1	3041129	04/11/13	04/12/13	RSK-175	
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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

LL_TB_041013
T130836-06 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041124	04/11/13	04/12/13	EPA 8260B
Bromochloromethane	ND	1.0	"	"	"	"	"	"
Bromodichloromethane	ND	1.0	"	"	"	"	"	"
Bromoform	ND	1.0	"	"	"	"	"	"
Bromomethane	ND	1.0	"	"	"	"	"	"
n-Butylbenzene	ND	1.0	"	"	"	"	"	"
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"
Chlorobenzene	ND	1.0	"	"	"	"	"	"
Chloroethane	ND	1.0	"	"	"	"	"	"
Chloroform	ND	1.0	"	"	"	"	"	"
Chloromethane	ND	1.0	"	"	"	"	"	"
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"
Dibromochloromethane	ND	1.0	"	"	"	"	"	"
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"
Dibromomethane	ND	1.0	"	"	"	"	"	"
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

LL_TB_041013
T130836-06 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

cis-1,3-Dichloropropene	ND	0.50	ug/l	1	3041124	04/11/13	04/12/13	EPA 8260B
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"
Tert-butyl alcohol	ND	10	"	"	"	"	"	"
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

LL_TB_041013
T130836-06 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	ug/l	1	3041124	04/11/13	04/12/13	EPA 8260B	
Surrogate: 4-Bromofluorobenzene	106 %	83.5-119			"	"	"	"	
Surrogate: Dibromofluoromethane	105 %	81-136			"	"	"	"	
Surrogate: Toluene-d8	100 %	88.8-117			"	"	"	"	

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Purgeable Petroleum Hydrocarbons by EPA 8015C - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
Batch 3041128 - EPA 5030 GC										
Blank (3041128-BLK1)				Prepared: 04/11/13 Analyzed: 04/12/13						
C6-C12 (GRO)	ND	50	ug/l							
Surrogate 4-Bromofluorobenzene	106		"	100		106	65-135			
LCS (3041128-BS1)				Prepared: 04/11/13 Analyzed: 04/12/13						
C6-C12 (GRO)	5210	50	ug/l	5520		94.4	75-125			
Surrogate 4-Bromofluorobenzene	77.2		"	100		77.2	65-135			
Matrix Spike (3041128-MS1)				Source: T130836-01		Prepared: 04/11/13 Analyzed: 04/12/13				
C6-C12 (GRO)	5900	50	ug/l	5520	1350	82.3	65-135			
Surrogate 4-Bromofluorobenzene	110		"	100		110	65-135			
Matrix Spike Dup (3041128-MSD1)				Source: T130836-01		Prepared: 04/11/13 Analyzed: 04/12/13				
C6-C12 (GRO)	5460	50	ug/l	5520	1350	74.5	65-135	7.61	20	
Surrogate 4-Bromofluorobenzene	103		"	100		103	65-135			

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

Metals by SM 3500 Series Methods - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041618 - EPA 3010A

Blank (3041618-BLK1)

Prepared & Analyzed: 04/16/13

Ferrous Iron	ND	0.100	mg/l
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Duplicate (3041618-DUP1)

Source: T130862-01

Prepared & Analyzed: 04/16/13

Ferrous Iron	0.0260	0.100	mg/l	0.0380	37.5	200
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SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041124 - EPA 5030 GCMS

Blank (3041124-BLK1)

Prepared & Analyzed: 04/11/13

Bromobenzene	ND	1.0	ug/l
Bromochloromethane	ND	1.0	"
Bromodichloromethane	ND	1.0	"
Bromoform	ND	1.0	"
Bromomethane	ND	1.0	"
n-Butylbenzene	ND	1.0	"
sec-Butylbenzene	ND	1.0	"
tert-Butylbenzene	ND	1.0	"
Carbon tetrachloride	ND	0.50	"
Chlorobenzene	ND	1.0	"
Chloroethane	ND	1.0	"
Chloroform	ND	1.0	"
Chloromethane	ND	1.0	"
2-Chlorotoluene	ND	1.0	"
4-Chlorotoluene	ND	1.0	"
Dibromochloromethane	ND	1.0	"
1,2-Dibromo-3-chloropropane	ND	1.0	"
1,2-Dibromoethane (EDB)	ND	1.0	"
Dibromomethane	ND	1.0	"
1,2-Dichlorobenzene	ND	1.0	"
1,3-Dichlorobenzene	ND	1.0	"
1,4-Dichlorobenzene	ND	1.0	"
Dichlorodifluoromethane	ND	0.50	"
1,1-Dichloroethane	ND	1.0	"
1,2-Dichloroethane	ND	0.50	"
1,1-Dichloroethene	ND	1.0	"
cis-1,2-Dichloroethene	ND	1.0	"
trans-1,2-Dichloroethene	ND	1.0	"
1,2-Dichloropropane	ND	1.0	"
1,3-Dichloropropane	ND	1.0	"
2,2-Dichloropropane	ND	1.0	"
1,1-Dichloropropene	ND	1.0	"
cis-1,3-Dichloropropene	ND	0.50	"
trans-1,3-Dichloropropene	ND	0.50	"
Hexachlorobutadiene	ND	1.0	"
Isopropylbenzene	ND	1.0	"

SunStar Laboratories, Inc.

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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041124 - EPA 5030 GCMS

Blank (3041124-BLK1)

Prepared & Analyzed: 04/11/13

p-Isopropyltoluene	ND	1.0	ug/l
Methylene chloride	ND	1.0	"
Naphthalene	ND	1.0	"
n-Propylbenzene	ND	1.0	"
Styrene	ND	1.0	"
1,1,2,2-Tetrachloroethane	ND	1.0	"
1,1,1,2-Tetrachloroethane	ND	1.0	"
Tetrachloroethene	ND	1.0	"
1,2,3-Trichlorobenzene	ND	1.0	"
1,2,4-Trichlorobenzene	ND	1.0	"
1,1,2-Trichloroethane	ND	1.0	"
1,1,1-Trichloroethane	ND	1.0	"
Trichloroethene	ND	1.0	"
Trichlorofluoromethane	ND	1.0	"
1,2,3-Trichloropropane	ND	1.0	"
1,3,5-Trimethylbenzene	ND	1.0	"
1,2,4-Trimethylbenzene	ND	1.0	"
Vinyl chloride	ND	1.0	"
Benzene	ND	0.50	"
Toluene	ND	0.50	"
Ethylbenzene	ND	0.50	"
m,p-Xylene	ND	1.0	"
o-Xylene	ND	0.50	"
Tert-amyl methyl ether	ND	2.0	"
Tert-butyl alcohol	ND	10	"
Di-isopropyl ether	ND	2.0	"
Ethyl tert-butyl ether	ND	2.0	"
Methyl tert-butyl ether	ND	1.0	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"
Surrogate 4-Bromofluorobenzene	8.47		"
Surrogate Dibromofluoromethane	8.13		"
Surrogate Toluene-d8	8.22		"

8.00 106 83.5-119
8.00 102 81-136
8.00 103 88.8-117

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041124 - EPA 5030 GCMS

LCS (3041124-BS1)

Prepared: 04/11/13 Analyzed: 04/12/13

Chlorobenzene	20.0	1.0	ug/l	20.0		100	75-125			
1,1-Dichloroethene	21.7	1.0	"	20.0		109	75-125			
Trichloroethene	20.8	1.0	"	20.0		104	75-125			
Benzene	20.5	0.50	"	20.0		102	75-125			
Toluene	19.3	0.50	"	20.0		96.4	75-125			
Surrogate 4-Bromofluorobenzene	8.35		"	8.00		104	83.5-119			
Surrogate Dibromofluoromethane	8.42		"	8.00		105	81-136			
Surrogate Toluene-d8	8.14		"	8.00		102	88.8-117			

Matrix Spike (3041124-MS1)

Source: T130836-01

Prepared: 04/11/13 Analyzed: 04/12/13

Chlorobenzene	19.6	1.0	ug/l	20.0	ND	98.1	75-125			
1,1-Dichloroethene	22.3	1.0	"	20.0	ND	112	75-125			
Trichloroethene	19.4	1.0	"	20.0	ND	96.8	75-125			
Benzene	20.5	0.50	"	20.0	ND	102	75-125			
Toluene	18.7	0.50	"	20.0	ND	93.6	75-125			
Surrogate 4-Bromofluorobenzene	7.42		"	8.00		92.8	83.5-119			
Surrogate Dibromofluoromethane	9.08		"	8.00		114	81-136			
Surrogate Toluene-d8	7.89		"	8.00		98.6	88.8-117			

Matrix Spike Dup (3041124-MSD1)

Source: T130836-01

Prepared: 04/11/13 Analyzed: 04/12/13

Chlorobenzene	20.0	1.0	ug/l	20.0	ND	100	75-125	2.07	20	
1,1-Dichloroethene	21.4	1.0	"	20.0	ND	107	75-125	4.03	20	
Trichloroethene	19.2	1.0	"	20.0	ND	95.9	75-125	0.882	20	
Benzene	20.5	0.50	"	20.0	ND	103	75-125	0.244	20	
Toluene	18.9	0.50	"	20.0	ND	94.4	75-125	0.957	20	
Surrogate 4-Bromofluorobenzene	7.52		"	8.00		94.0	83.5-119			
Surrogate Dibromofluoromethane	8.52		"	8.00		106	81-136			
Surrogate Toluene-d8	8.00		"	8.00		100	88.8-117			

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/17/13 14:10
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Conventional Chemistry Parameters by APHA/EPA/ASTM Methods - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041122 - General Preparation

Duplicate (3041122-DUP1) **Source: T130836-02** Prepared: 04/11/13 Analyzed: 04/12/13

Total Alkalinity	570	20	mg/l		555			2.67	25	
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SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/17/13 14:10
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Anions by EPA Method 300.0 - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
Batch 3041121 - General Preparation										
Blank (3041121-BLK1)				Prepared & Analyzed: 04/11/13						
Sulfate as SO4	ND	0.500	mg/l							
Nitrate as NO3	ND	0.500	"							
LCS (3041121-BS1)				Prepared & Analyzed: 04/11/13						
Sulfate as SO4	10.2	0.500	mg/l	10.0		102	75-125			
Nitrate as NO3	0.604	0.500	"	0.500		121	75-125			
Matrix Spike (3041121-MS1)				Source: T130836-02 Prepared & Analyzed: 04/11/13						
Sulfate as SO4	379	0.500	mg/l	10.0	383	NR	75-125			QM-02
Nitrate as NO3	0.621	0.500	"	0.500	ND	124	75-125			
Matrix Spike Dup (3041121-MSD1)				Source: T130836-02 Prepared & Analyzed: 04/11/13						
Sulfate as SO4	379	0.500	mg/l	10.0	383	NR	75-125	0.0314	20	QM-02
Nitrate as NO3	0.617	0.500	"	0.500	ND	123	75-125	0.646	20	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

RSK-175 - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041129 - EPA 3810m Headspace

Blank (3041129-BLK1)

Prepared: 04/11/13 Analyzed: 04/12/13

Methane	ND	1.00	ug/l							
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Duplicate (3041129-DUP1)

Source: T130836-02

Prepared: 04/11/13 Analyzed: 04/12/13

Methane	1.19	1.00	ug/l		1.17			1.69	20	
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SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 14:10

Notes and Definitions

- S-04 The surrogate recovery for this sample is outside of established control limits due to a sample matrix effect.
- QM-02 The RPD and/or percent recovery for this QC spike sample cannot be accurately calculated due to the high concentration of analyte inherent in the sample.
- DET Analyte DETECTED
- ND Analyte NOT DETECTED at or above the reporting limit
- NR Not Reported
- dry Sample results reported on a dry weight basis
- RPD Relative Percent Difference

SunStar Laboratories, Inc.

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Katherine Shields

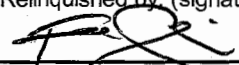

Katherine Shields, Jr. Project Manager

SunStar Laboratories, Inc.
25712 Commercentre Dr
Lake Forest, CA 92630
949-297-5020

Chain of Custody Record

Client: MUREX ENVIRONMENTAL INC.
Address: 2640 Walnut Ave, Unit F
Phone: (714) 508-0800 Fax: (714) 508-0880
Project Manager: Jeremy Squire (714) 604-5836

Date: 4-10-13 Page: 1 OF 1
Project Name: CENCO
Collector: Frane Sosic +AW Client Project #: 1003-001-300
Batch #: T130836 EDF #:

Sample ID	Date Sampled	Time	Sample Type	TPHg (8015 M)	VOCs (8260 B)	Methane (8015)	Alkalinity	300 IC (Nitrate, Sulfate)	Ferrous Iron									Total # of containers	Comments/Preservative	Laboratory ID #
LL-EWL-041013	4-10-13	1000	GW	X	X													6		01
LL-703-041013	4-10-13	1245	GW	X	X	X	X	X	X									12		02
LL-701-041013	4-10-13	1413	GW	X	X													6		03
LL-702-041013	4-10-13	1542	GW	X	X	X	X	X	X									12		04
LL-705-041013	4-10-13	1636	GW	X	X	X	X	X	X									12		05
LL-TB-041013			Water															2		06
Relinquished by: (signature) 	Date / Time 4-10-13 17:00	Received by: (Sign / Date / Time)  4-10-13 17:00	Total # of containers 50	Notes 3.4																
Relinquished by: (signature)	Date / Time	Received by: (Sign / Date / Time)	Chain of Custody seals N																	
Relinquished by: (signature)	Date / Time	Received by: (Sign / Date / Time)	Seals intact? Y/N/NA N/A																	
Relinquished by: (signature)	Date / Time	Received by: (Sign / Date / Time)	Received good condition/cold Y																	
			Turn around time: Standard																	

Sample disposal Instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____

SAMPLE RECEIVING REVIEW SHEET

BATCH # T130836

Client Name: Murex

Project: Cenco

Received by: Patrick

Date/Time Received: 4/10/13 1700

Delivered by: ☐ Client ☒ SunStar Courier ☐ GSO ☐ FedEx ☐ Other _____

Total number of coolers received 1 Temp criteria = 6°C > 0°C (no frozen containers)

Temperature: cooler #1 3.6 °C +/- the CF (- 0.2°C) = 3.4 °C corrected temperature

cooler #2 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

cooler #3 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

Samples outside temp. but received on ice, w/in 6 hours of final sampling. ☒ Yes ☐ No* ☐ N/A

Custody Seals Intact on Cooler/Sample ☐ Yes ☐ No* ☒ N/A

Sample Containers Intact ☒ Yes ☐ No*

Sample labels match COC ID's ☒ Yes ☐ No*


Total number of containers received match COC ☒ Yes ☐ No*

Proper containers received for analyses requested on COC ☒ Yes ☐ No*

Proper preservative indicated on COC/containers for analyses requested ☒ Yes ☐ No* ☐ N/A

Complete shipment received in good condition with correct temperatures, containers, labels, volumes preservatives and within method specified holding times. ☒ Yes ☐ No*

* Complete Non-Conformance Receiving Sheet if checked

Cooler/Sample Review - Initials and date  4/10/13

Comments:



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17 April 2013

Jeremy Squire
Murex
15375 Barranca Parkway, Suite K-101
Irvine, CA 92861
RE: Cenco

Enclosed are the results of analyses for samples received by the laboratory on 04/11/13 17:00. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Katherine Shields
Jr. Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
LL_714_041113	T130847-01	Water	04/11/13 10:05	04/11/13 17:00
LL_713_041113_01	T130847-02	Water	04/11/13 10:05	04/11/13 17:00
LL_713_041113_02	T130847-03	Water	04/11/13 11:05	04/11/13 17:00
LL_709_041113	T130847-04	Water	04/11/13 13:50	04/11/13 17:00
LL_708_041113_01	T130847-05	Water	04/11/13 16:15	04/11/13 17:00
LL_708_041113_02	T130847-06	Water	04/11/13 16:20	04/11/13 17:00
LL_TB_041113	T130847-07	Water	04/11/13 00:00	04/11/13 17:00

SunStar Laboratories, Inc.

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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_714_041113
T130847-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	170	50	ug/l	1	3041214	04/12/13	04/15/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		108 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041208	04/12/13	04/13/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	2.7	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager

Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_714_041113
T130847-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3041208	04/12/13	04/13/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	7.8	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	1.2	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	1.3	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_714_041113
T130847-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	52	10	ug/l	1	3041208	04/12/13	04/13/13	EPA 8260B	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	7.2	1.0	"	"	"	"	"	"	
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene		109 %	83.5-119		"	"	"	"	
Surrogate: Dibromofluoromethane		107 %	81-136		"	"	"	"	
Surrogate: Toluene-d8		104 %	88.8-117		"	"	"	"	

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_713_041113_01
T130847-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	8200	50	ug/l	1	3041214	04/12/13	04/15/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		82.0 %	65-135		"	"	"	"	

Metals by SM 3500 Series Methods

Ferrous Iron	0.586	0.100	mg/l	1	3041618	04/15/13	04/16/13	EPA6010/S M3500	
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Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041208	04/12/13	04/13/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	11	1.0	"	"	"	"	"	"	
sec-Butylbenzene	15	1.0	"	"	"	"	"	"	
tert-Butylbenzene	1.6	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_713_041113_01
T130847-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1-Dichloroethene	ND	1.0	ug/l	1	3041208	04/12/13	04/13/13	EPA 8260B	
cis-1,2-Dichloroethene	2.0	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	110	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	6.8	1.0	"	"	"	"	"	"	
n-Propylbenzene	190	1.0	"	"	"	"	"	"	E-1
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	13	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	1.5	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	4900	50	"	100	"	"	"	"	
Toluene	8.2	0.50	"	1	"	"	"	"	
Ethylbenzene	13	0.50	"	"	"	"	"	"	
m,p-Xylene	37	1.0	"	"	"	"	"	"	

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Katherine Shields

Katherine Shields, Jr. Project Manager

Murex
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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_713_041113_01
T130847-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

o-Xylene	1.9	0.50	ug/l	1	3041208	04/12/13	04/13/13	EPA 8260B	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	310	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	760	1.0	"	"	"	"	"	"	E-1
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene		105 %	83.5-119		"	"	"	"	
Surrogate: Dibromofluoromethane		105 %	81-136		"	"	"	"	
Surrogate: Toluene-d8		102 %	88.8-117		"	"	"	"	

Conventional Chemistry Parameters by APHA/EPA/ASTM Methods

Total Alkalinity	840	20	mg/l	1	3041224	04/12/13	04/12/13	EPA 310.1	
Anions by EPA Method 300.0									
Sulfate as SO4	19.7	0.500	mg/l	1	3041204	04/12/13	04/12/13	EPA 300.0	
Nitrate as NO3	ND	0.500	"	"	"	"	"	"	
RSK-175									
Methane	4220	10.0	ug/l	10	3041211	04/12/13	04/12/13	RSK-175	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_713_041113_02
T130847-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	8300	50	ug/l	1	3041214	04/12/13	04/15/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		80.0 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041208	04/12/13	04/13/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	11	1.0	"	"	"	"	"	"	
sec-Butylbenzene	15	1.0	"	"	"	"	"	"	
tert-Butylbenzene	1.6	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	2.1	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_713_041113_02
T130847-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,3-Dichloropropane	ND	1.0	ug/l	1	3041208	04/12/13	04/13/13	EPA 8260B	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	110	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	6.7	1.0	"	"	"	"	"	"	
n-Propylbenzene	190	1.0	"	"	"	"	"	"	E-1
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	14	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	1.5	1.0	"	"	"	"	"	"	
Vinyl chloride	1.2	1.0	"	"	"	"	"	"	
Benzene	5000	50	"	100	"	"	"	"	
Toluene	8.4	0.50	"	1	"	"	"	"	
Ethylbenzene	13	0.50	"	"	"	"	"	"	
m,p-Xylene	38	1.0	"	"	"	"	"	"	
o-Xylene	2.0	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	320	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_713_041113_02
T130847-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Ethyl tert-butyl ether	ND	2.0	ug/l	1	3041208	04/12/13	04/13/13	EPA 8260B	
Methyl tert-butyl ether	800	1.0	"	"	"	"	"	"	E-1
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene		108 %	83.5-119		"	"	"	"	
Surrogate: Dibromofluoromethane		108 %	81-136		"	"	"	"	
Surrogate: Toluene-d8		101 %	88.8-117		"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_709_041113
T130847-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	850	50	ug/l	1	3041214	04/12/13	04/15/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		117 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041208	04/12/13	04/15/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	4.9	1.0	"	"	"	"	"	"	
tert-Butylbenzene	1.0	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_709_041113
T130847-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,3-Dichloropropane	ND	1.0	ug/l	1	3041208	04/12/13	04/15/13	EPA 8260B
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	25	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	20	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"
Tert-butyl alcohol	160	10	"	"	"	"	"	"

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Page 12 of 34



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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_709_041113
T130847-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Di-isopropyl ether	ND	2.0	ug/l	1	3041208	04/12/13	04/15/13	EPA 8260B
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	2.1	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		104 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		92.4 %	81-136		"	"	"	"
Surrogate: Toluene-d8		103 %	88.8-117		"	"	"	"

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/17/13 17:00
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LL_708_041113_01
T130847-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	25000	250	ug/l	5	3041214	04/12/13	04/15/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		91.9 %	65-135		"	"	"	"	

Metals by SM 3500 Series Methods

Ferrous Iron	1.23	0.100	mg/l	1	3041618	04/15/13	04/16/13	EPA6010/S M3500	
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Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041208	04/12/13	04/13/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	7.6	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	2.5	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_708_041113_01
T130847-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1-Dichloroethene	ND	1.0	ug/l	1	3041208	04/12/13	04/13/13	EPA 8260B	
cis-1,2-Dichloroethene	2.4	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	18	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	27	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	290	1.0	"	"	"	"	"	"	E-1
n-Propylbenzene	9.4	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	670	50	"	50	"	"	"	"	
1,2,4-Trimethylbenzene	1700	50	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	1	"	"	"	"	
Benzene	1100	25	"	50	"	"	"	"	
Toluene	54	0.50	"	1	"	"	"	"	
Ethylbenzene	510	25	"	50	"	"	"	"	
m,p-Xylene	920	50	"	"	"	"	"	"	

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Katherine Shields

Katherine Shields, Jr. Project Manager

Murex
15375 Barranca Parkway, Suite K-101
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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_708_041113_01
T130847-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

o-Xylene	27	0.50	ug/l	1	3041208	04/12/13	04/13/13	EPA 8260B	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	350	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	790	1.0	"	"	"	"	"	"	E-1
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene		80.9 %	83.5-119		"	"	"	"	S-GC
Surrogate: Dibromofluoromethane		102 %	81-136		"	"	"	"	
Surrogate: Toluene-d8		92.2 %	88.8-117		"	"	"	"	

Conventional Chemistry Parameters by APHA/EPA/ASTM Methods

Total Alkalinity	860	20	mg/l	1	3041224	04/12/13	04/12/13	EPA 310.1	
Anions by EPA Method 300.0									
Sulfate as SO4	48.4	0.500	mg/l	1	3041204	04/12/13	04/12/13	EPA 300.0	
Nitrate as NO3	ND	0.500	"	"	"	"	"	"	
RSK-175									
Methane	4110	10.0	ug/l	10	3041211	04/12/13	04/12/13	RSK-175	

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Project: Cenco
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Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_708_041113_02
T130847-06 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	240000	2500	ug/l	50	3041214	04/12/13	04/15/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		146 %	65-135		"	"	"	"	S-01

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041208	04/12/13	04/15/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	130	1.0	"	"	"	"	"	"	E-1
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	2.4	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_708_041113_02
T130847-06 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,3-Dichloropropane	ND	1.0	ug/l	1	3041208	04/12/13	04/15/13	EPA 8260B	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	65	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	120	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	680	1.0	"	"	"	"	"	"	E-1
n-Propylbenzene	140	50	"	50	"	"	"	"	
Styrene	ND	1.0	"	1	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	780	50	"	50	"	"	"	"	
1,2,4-Trimethylbenzene	2000	50	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	1	"	"	"	"	
Benzene	990	25	"	50	"	"	"	"	
Toluene	54	0.50	"	1	"	"	"	"	
Ethylbenzene	430	25	"	50	"	"	"	"	
m,p-Xylene	890	50	"	"	"	"	"	"	
o-Xylene	24	0.50	"	1	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	260	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	

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Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_708_041113_02
T130847-06 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Ethyl tert-butyl ether	ND	2.0	ug/l	1	3041208	04/12/13	04/15/13	EPA 8260B	
Methyl tert-butyl ether	670	1.0	"	"	"	"	"	"	E-1
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene		42.8 %	83.5-119		"	"	"	"	S-GC
Surrogate: Dibromofluoromethane		94.8 %	81-136		"	"	"	"	
Surrogate: Toluene-d8		95.5 %	88.8-117		"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_TB_041113
T130847-07 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041526	04/15/13	04/16/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	

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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_TB_041113
T130847-07 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

cis-1,3-Dichloropropene	ND	0.50	ug/l	1	3041526	04/15/13	04/16/13	EPA 8260B
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"
Tert-butyl alcohol	ND	10	"	"	"	"	"	"
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"

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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

LL_TB_041113
T130847-07 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	ug/l	1	3041526	04/15/13	04/16/13	EPA 8260B	
Surrogate: 4-Bromofluorobenzene	107 %	83.5-119			"	"	"	"	
Surrogate: Dibromofluoromethane	96.6 %	81-136			"	"	"	"	
Surrogate: Toluene-d8	105 %	88.8-117			"	"	"	"	

SunStar Laboratories, Inc.

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/17/13 17:00
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Purgeable Petroleum Hydrocarbons by EPA 8015C - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
Batch 3041214 - EPA 5030 GC										
Blank (3041214-BLK1)				Prepared: 04/12/13 Analyzed: 04/15/13						
C6-C12 (GRO)	ND	50	ug/l							
Surrogate 4-Bromofluorobenzene	121		"	100		121	65-135			
LCS (3041214-BS1)				Prepared: 04/12/13 Analyzed: 04/15/13						
C6-C12 (GRO)	5000	50	ug/l	5520		90.6	75-125			
Surrogate 4-Bromofluorobenzene	93.7		"	100		93.7	65-135			
Matrix Spike (3041214-MS1)				Source: T130851-03		Prepared: 04/12/13 Analyzed: 04/15/13				
C6-C12 (GRO)	4610	50	ug/l	5520	ND	83.6	65-135			
Surrogate 4-Bromofluorobenzene	88.4		"	100		88.4	65-135			
Matrix Spike Dup (3041214-MSD1)				Source: T130851-03		Prepared: 04/12/13 Analyzed: 04/15/13				
C6-C12 (GRO)	4780	50	ug/l	5520	ND	86.6	65-135	3.53	20	
Surrogate 4-Bromofluorobenzene	103		"	100		103	65-135			

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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

Metals by SM 3500 Series Methods - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041618 - EPA 3010A

Blank (3041618-BLK1)

Prepared & Analyzed: 04/16/13

Ferrous Iron	ND	0.100	mg/l
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Duplicate (3041618-DUP1)

Source: T130862-01

Prepared & Analyzed: 04/16/13

Ferrous Iron	0.0260	0.100	mg/l	0.0380	37.5	200
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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041208 - EPA 5030 GCMS

Blank (3041208-BLK1)

Prepared: 04/12/13 Analyzed: 04/13/13

Bromobenzene	ND	1.0	ug/l
Bromochloromethane	ND	1.0	"
Bromodichloromethane	ND	1.0	"
Bromoform	ND	1.0	"
Bromomethane	ND	1.0	"
n-Butylbenzene	ND	1.0	"
sec-Butylbenzene	ND	1.0	"
tert-Butylbenzene	ND	1.0	"
Carbon tetrachloride	ND	0.50	"
Chlorobenzene	ND	1.0	"
Chloroethane	ND	1.0	"
Chloroform	ND	1.0	"
Chloromethane	ND	1.0	"
2-Chlorotoluene	ND	1.0	"
4-Chlorotoluene	ND	1.0	"
Dibromochloromethane	ND	1.0	"
1,2-Dibromo-3-chloropropane	ND	1.0	"
1,2-Dibromoethane (EDB)	ND	1.0	"
Dibromomethane	ND	1.0	"
1,2-Dichlorobenzene	ND	1.0	"
1,3-Dichlorobenzene	ND	1.0	"
1,4-Dichlorobenzene	ND	1.0	"
Dichlorodifluoromethane	ND	0.50	"
1,1-Dichloroethane	ND	1.0	"
1,2-Dichloroethane	ND	0.50	"
1,1-Dichloroethene	ND	1.0	"
cis-1,2-Dichloroethene	ND	1.0	"
trans-1,2-Dichloroethene	ND	1.0	"
1,2-Dichloropropane	ND	1.0	"
1,3-Dichloropropane	ND	1.0	"
2,2-Dichloropropane	ND	1.0	"
1,1-Dichloropropene	ND	1.0	"
cis-1,3-Dichloropropene	ND	0.50	"
trans-1,3-Dichloropropene	ND	0.50	"
Hexachlorobutadiene	ND	1.0	"
Isopropylbenzene	ND	1.0	"

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/17/13 17:00
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Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041208 - EPA 5030 GCMS

Blank (3041208-BLK1)

Prepared: 04/12/13 Analyzed: 04/13/13

p-Isopropyltoluene	ND	1.0	ug/l							
Methylene chloride	ND	1.0	"							
Naphthalene	ND	1.0	"							
n-Propylbenzene	ND	1.0	"							
Styrene	ND	1.0	"							
1,1,2,2-Tetrachloroethane	ND	1.0	"							
1,1,1,2-Tetrachloroethane	ND	1.0	"							
Tetrachloroethene	ND	1.0	"							
1,2,3-Trichlorobenzene	ND	1.0	"							
1,2,4-Trichlorobenzene	ND	1.0	"							
1,1,2-Trichloroethane	ND	1.0	"							
1,1,1-Trichloroethane	ND	1.0	"							
Trichloroethene	ND	1.0	"							
Trichlorofluoromethane	ND	1.0	"							
1,2,3-Trichloropropane	ND	1.0	"							
1,3,5-Trimethylbenzene	ND	1.0	"							
1,2,4-Trimethylbenzene	ND	1.0	"							
Vinyl chloride	ND	1.0	"							
Benzene	ND	0.50	"							
Toluene	ND	0.50	"							
Ethylbenzene	ND	0.50	"							
m,p-Xylene	ND	1.0	"							
o-Xylene	ND	0.50	"							
Tert-amyl methyl ether	ND	2.0	"							
Tert-butyl alcohol	ND	10	"							
Di-isopropyl ether	ND	2.0	"							
Ethyl tert-butyl ether	ND	2.0	"							
Methyl tert-butyl ether	ND	1.0	"							
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"							
Surrogate 4-Bromofluorobenzene	8.72		"	8.00		109	83.5-119			
Surrogate Dibromofluoromethane	8.53		"	8.00		107	81-136			
Surrogate Toluene-d8	8.23		"	8.00		103	88.8-117			

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/17/13 17:00
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Volatile Organic Compounds by EPA Method 8260B - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041208 - EPA 5030 GCMS

LCS (3041208-BS1)

Prepared: 04/12/13 Analyzed: 04/13/13

Chlorobenzene	20.1	1.0	ug/l	20.0		100	75-125			
1,1-Dichloroethene	21.9	1.0	"	20.0		110	75-125			
Trichloroethene	20.9	1.0	"	20.0		104	75-125			
Benzene	21.9	0.50	"	20.0		110	75-125			
Toluene	17.3	0.50	"	20.0		86.4	75-125			
Surrogate 4-Bromofluorobenzene	8.40		"	8.00		105	83.5-119			
Surrogate Dibromofluoromethane	9.59		"	8.00		120	81-136			
Surrogate Toluene-d8	7.80		"	8.00		97.5	88.8-117			

Matrix Spike (3041208-MS1)

Source: T130847-01

Prepared: 04/12/13 Analyzed: 04/13/13

Chlorobenzene	19.3	1.0	ug/l	20.0	ND	96.7	75-125			
1,1-Dichloroethene	20.3	1.0	"	20.0	ND	101	75-125			
Trichloroethene	18.7	1.0	"	20.0	ND	93.3	75-125			
Benzene	22.6	0.50	"	20.0	1.34	106	75-125			
Toluene	18.5	0.50	"	20.0	ND	92.4	75-125			
Surrogate 4-Bromofluorobenzene	8.44		"	8.00		106	83.5-119			
Surrogate Dibromofluoromethane	9.21		"	8.00		115	81-136			
Surrogate Toluene-d8	8.24		"	8.00		103	88.8-117			

Matrix Spike Dup (3041208-MSD1)

Source: T130847-01

Prepared: 04/12/13 Analyzed: 04/13/13

Chlorobenzene	19.6	1.0	ug/l	20.0	ND	98.2	75-125	1.59	20	
1,1-Dichloroethene	19.5	1.0	"	20.0	ND	97.7	75-125	3.77	20	
Trichloroethene	18.8	1.0	"	20.0	ND	94.2	75-125	0.907	20	
Benzene	22.4	0.50	"	20.0	1.34	106	75-125	0.533	20	
Toluene	18.9	0.50	"	20.0	ND	94.6	75-125	2.35	20	
Surrogate 4-Bromofluorobenzene	8.33		"	8.00		104	83.5-119			
Surrogate Dibromofluoromethane	9.28		"	8.00		116	81-136			
Surrogate Toluene-d8	8.41		"	8.00		105	88.8-117			

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Murex	Project: Cenco	
15375 Barranca Parkway, Suite K-101	Project Number: 1003-001-300	
Irvine CA, 92861	Project Manager: Jeremy Squire	Reported: 04/17/13 17:00

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041526 - EPA 5030 GCMS

Blank (3041526-BLK1)

Prepared: 04/15/13 Analyzed: 04/16/13

Bromobenzene	ND	1.0	ug/l
Bromochloromethane	ND	1.0	"
Bromodichloromethane	ND	1.0	"
Bromoform	ND	1.0	"
Bromomethane	ND	1.0	"
n-Butylbenzene	ND	1.0	"
sec-Butylbenzene	ND	1.0	"
tert-Butylbenzene	ND	1.0	"
Carbon tetrachloride	ND	0.50	"
Chlorobenzene	ND	1.0	"
Chloroethane	ND	1.0	"
Chloroform	ND	1.0	"
Chloromethane	ND	1.0	"
2-Chlorotoluene	ND	1.0	"
4-Chlorotoluene	ND	1.0	"
Dibromochloromethane	ND	1.0	"
1,2-Dibromo-3-chloropropane	ND	1.0	"
1,2-Dibromoethane (EDB)	ND	1.0	"
Dibromomethane	ND	1.0	"
1,2-Dichlorobenzene	ND	1.0	"
1,3-Dichlorobenzene	ND	1.0	"
1,4-Dichlorobenzene	ND	1.0	"
Dichlorodifluoromethane	ND	0.50	"
1,1-Dichloroethane	ND	1.0	"
1,2-Dichloroethane	ND	0.50	"
1,1-Dichloroethene	ND	1.0	"
cis-1,2-Dichloroethene	ND	1.0	"
trans-1,2-Dichloroethene	ND	1.0	"
1,2-Dichloropropane	ND	1.0	"
1,3-Dichloropropane	ND	1.0	"
2,2-Dichloropropane	ND	1.0	"
1,1-Dichloropropene	ND	1.0	"
cis-1,3-Dichloropropene	ND	0.50	"
trans-1,3-Dichloropropene	ND	0.50	"
Hexachlorobutadiene	ND	1.0	"
Isopropylbenzene	ND	1.0	"

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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041526 - EPA 5030 GCMS

Blank (3041526-BLK1)

Prepared: 04/15/13 Analyzed: 04/16/13

p-Isopropyltoluene	ND	1.0	ug/l
Methylene chloride	ND	1.0	"
Naphthalene	ND	1.0	"
n-Propylbenzene	ND	1.0	"
Styrene	ND	1.0	"
1,1,2,2-Tetrachloroethane	ND	1.0	"
1,1,1,2-Tetrachloroethane	ND	1.0	"
Tetrachloroethene	ND	1.0	"
1,2,3-Trichlorobenzene	ND	1.0	"
1,2,4-Trichlorobenzene	ND	1.0	"
1,1,2-Trichloroethane	ND	1.0	"
1,1,1-Trichloroethane	ND	1.0	"
Trichloroethene	ND	1.0	"
Trichlorofluoromethane	ND	1.0	"
1,2,3-Trichloropropane	ND	1.0	"
1,3,5-Trimethylbenzene	ND	1.0	"
1,2,4-Trimethylbenzene	ND	1.0	"
Vinyl chloride	ND	1.0	"
Benzene	ND	0.50	"
Toluene	ND	0.50	"
Ethylbenzene	ND	0.50	"
m,p-Xylene	ND	1.0	"
o-Xylene	ND	0.50	"
Tert-amyl methyl ether	ND	2.0	"
Tert-butyl alcohol	ND	10	"
Di-isopropyl ether	ND	2.0	"
Ethyl tert-butyl ether	ND	2.0	"
Methyl tert-butyl ether	ND	1.0	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"
Surrogate 4-Bromofluorobenzene	8.26		"
Surrogate Dibromofluoromethane	7.73		"
Surrogate Toluene-d8	8.51		"

8.00 103 83.5-119
8.00 96.6 81-136
8.00 106 88.8-117

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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041526 - EPA 5030 GCMS

LCS (3041526-BS1)

Prepared: 04/15/13 Analyzed: 04/16/13

Chlorobenzene	20.0	1.0	ug/l	20.0		99.8	75-125			
1,1-Dichloroethene	22.2	1.0	"	20.0		111	75-125			
Trichloroethene	20.8	1.0	"	20.0		104	75-125			
Benzene	21.1	0.50	"	20.0		106	75-125			
Toluene	19.7	0.50	"	20.0		98.7	75-125			
Surrogate 4-Bromofluorobenzene	8.47		"	8.00		106	83.5-119			
Surrogate Dibromofluoromethane	8.25		"	8.00		103	81-136			
Surrogate Toluene-d8	8.32		"	8.00		104	88.8-117			

Matrix Spike (3041526-MS1)

Source: T130856-03

Prepared: 04/15/13 Analyzed: 04/16/13

Chlorobenzene	19.6	1.0	ug/l	20.0	ND	98.0	75-125			
1,1-Dichloroethene	22.5	1.0	"	20.0	ND	113	75-125			
Trichloroethene	20.2	1.0	"	20.0	ND	101	75-125			
Benzene	21.5	0.50	"	20.0	ND	108	75-125			
Toluene	19.8	0.50	"	20.0	ND	99.1	75-125			
Surrogate 4-Bromofluorobenzene	8.37		"	8.00		105	83.5-119			
Surrogate Dibromofluoromethane	8.67		"	8.00		108	81-136			
Surrogate Toluene-d8	8.29		"	8.00		104	88.8-117			

Matrix Spike Dup (3041526-MSD1)

Source: T130856-03

Prepared: 04/15/13 Analyzed: 04/16/13

Chlorobenzene	19.4	1.0	ug/l	20.0	ND	97.2	75-125	0.717	20	
1,1-Dichloroethene	22.2	1.0	"	20.0	ND	111	75-125	1.38	20	
Trichloroethene	20.3	1.0	"	20.0	ND	102	75-125	0.444	20	
Benzene	21.4	0.50	"	20.0	ND	107	75-125	0.233	20	
Toluene	19.9	0.50	"	20.0	ND	99.6	75-125	0.553	20	
Surrogate 4-Bromofluorobenzene	8.55		"	8.00		107	83.5-119			
Surrogate Dibromofluoromethane	8.55		"	8.00		107	81-136			
Surrogate Toluene-d8	8.40		"	8.00		105	88.8-117			

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/17/13 17:00
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Conventional Chemistry Parameters by APHA/EPA/ASTM Methods - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041224 - General Preparation

Duplicate (3041224-DUP1) **Source: T130847-05** Prepared & Analyzed: 04/12/13

Total Alkalinity	850	20	mg/l		855			0.587	25	
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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/17/13 17:00
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Anions by EPA Method 300.0 - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041204 - General Preparation

Blank (3041204-BLK1)				Prepared: 04/12/13 Analyzed: 04/15/13						
Fluoride	ND	0.500	mg/l							
Chloride	ND	0.500	"							
Nitrite as NO2	ND	0.500	"							
Sulfate as SO4	ND	0.500	"							
Nitrate as NO3	ND	0.500	"							

LCS (3041204-BS1)				Prepared & Analyzed: 04/12/13						
Chloride	10.3	0.500	mg/l	10.0		103	75-125			
Sulfate as SO4	10.8	0.500	"	10.0		108	75-125			
Nitrate as NO3	0.579	0.500	"	0.500		116	75-125			

Matrix Spike (3041204-MS1)				Source: T130845-11		Prepared & Analyzed: 04/12/13				
Chloride	17.9	0.500	mg/l	10.0	7.31	106	75-125			
Sulfate as SO4	10.3	0.500	"	10.0	0.962	93.0	75-125			
Nitrate as NO3	0.732	0.500	"	0.500	0.247	97.0	75-125			

Matrix Spike Dup (3041204-MSD1)				Source: T130845-11		Prepared & Analyzed: 04/12/13				
Chloride	17.8	0.500	mg/l	10.0	7.31	105	75-125	0.527	20	
Sulfate as SO4	10.2	0.500	"	10.0	0.962	92.7	75-125	0.371	20	
Nitrate as NO3	0.728	0.500	"	0.500	0.247	96.2	75-125	0.548	20	

SunStar Laboratories, Inc.

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15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/17/13 17:00

RSK-175 - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041211 - EPA 3810m Headspace

Blank (3041211-BLK1)

Prepared & Analyzed: 04/12/13

Methane	ND	1.00	ug/l
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Duplicate (3041211-DUP1)

Source: T130847-02

Prepared & Analyzed: 04/12/13

Methane	4190	10.0	ug/l	4220	0.652	20
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Katherine Shields, Jr. Project Manager



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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/17/13 17:00
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Notes and Definitions

S-GC Surrogate recovery outside of established control limits. The data was accepted based on valid recovery of the remaining surrogate(s).

S-01 The surrogate recovery for this sample is not available due to sample dilution required from high analyte concentration and/or matrix interference's.

E-1 The final dilution was lower than the original data or previous dilutions. The highest recovered concentration was reported even though it was above calibration range.

DET Analyte DETECTED

ND Analyte NOT DETECTED at or above the reporting limit

NR Not Reported

dry Sample results reported on a dry weight basis

RPD Relative Percent Difference

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager

SunStar Laboratories, Inc.
25712 Commercentre Dr
Lake Forest, CA 92630
949-297-5020

Chain of Custody Record

Client: MUREX ENVIRONMENTAL INC.

Address: 2640 Walnut Ave, Unit F

Phone: (714) 508-0800 Fax: (714) 508-0880

Project Manager: Jeremy Squire (714) 604-5836

Date: 04/11/2013

Project Name: CENCO

Collector: ~~Frank Sosie~~ AW

Batch #: 7130847

Page: 1 OF 1

Client Project #: 1003-001-300

EDF #: _____

Sample ID	Date Sampled	Time	Sample Type	TPHg (8015 M)	VOCs (8260 B)	METHANE (8015)	ALKALINITY	3001C (NITRATE, SULFATE)	FERRIC IRON									Total # of containers	Comments/Preservative	Laboratory ID #
LL-714-041113	04/11/13	1005	GW	X	X													6		01
LL-713-041113-01		1055		X	X	X	X	X	X									12		02
LL-713-041113-02		1105		X	X													6		03
LL-709-041113		1350		X	X													6		04
LL-708-041113-01		1615		X	X	X	X	X	X									12		05
LL-708-041113-02	↓	1620	↓	X	X													6		06
LL-TB-041113	-	-	WATER															2		07
Relinquished by: (signature) <i>T.A.L.</i>	Date / Time 04/11/13 1700	Received by: (Sign / Date / Time) <i>[Signature]</i> 4/11/13 1700	Total # of containers: 50																Notes	
Relinquished by: (signature)	Date / Time	Received by: (Sign / Date / Time)	Chain of Custody seals: NA																	
Relinquished by: (signature)	Date / Time	Received by: (Sign / Date / Time)	Seals intact? Y/N/NA: NA																	
Relinquished by: (signature)	Date / Time	Received by: (Sign / Date / Time)	Received good condition/cold: 3.2																	
			Turn around time: Standard																	

Sample disposal Instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____

SAMPLE RECEIVING REVIEW SHEET

BATCH # T130397

Client Name: MUREX

Project: CENCO

Received by: PATRICK

Date/Time Received: 4/11/13 17:00

Delivered by: ☐ Client ☒ SunStar Courier ☐ GSO ☐ FedEx ☐ Other _____

Total number of coolers received 0 Temp criteria = $6^{\circ}\text{C} > 0^{\circ}\text{C}$ (no frozen containers)

Temperature: cooler #1 3.4 $^{\circ}\text{C}$ +/- the CF (- 0.2 $^{\circ}\text{C}$) = 3.2 $^{\circ}\text{C}$ corrected temperature

cooler #2 _____ $^{\circ}\text{C}$ +/- the CF (- 0.2 $^{\circ}\text{C}$) = _____ $^{\circ}\text{C}$ corrected temperature

cooler #3 _____ $^{\circ}\text{C}$ +/- the CF (- 0.2 $^{\circ}\text{C}$) = _____ $^{\circ}\text{C}$ corrected temperature

Samples outside temp. but received on ice, w/in 6 hours of final sampling. ☒ Yes ☐ No* ☐ N/A

Custody Seals Intact on Cooler/Sample ☐ Yes ☐ No* ☒ N/A

Sample Containers Intact ☒ Yes ☐ No*

Sample labels match COC ID's ☒ Yes ☐ No*

Total number of containers received match COC ☒ Yes ☐ No*

Proper containers received for analyses requested on COC ☒ Yes ☐ No*

Proper preservative indicated on COC/containers for analyses requested ☒ Yes ☐ No* ☐ N/A

Complete shipment received in good condition with correct temperatures, containers, labels, volumes preservatives and within method specified holding times. ☒ Yes ☐ No*

* Complete Non-Conformance Receiving Sheet if checked

Cooler/Sample Review - Initials and date BC 4/12/13

Comments:



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19 April 2013

Jeremy Squire
Murex
15375 Barranca Parkway, Suite K-101
Irvine, CA 92861
RE: Cenco

Enclosed are the results of analyses for samples received by the laboratory on 04/12/13 16:34. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Wendy Hsiao For Katherine Shields
Jr. Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
LL_710_041213	T130856-01	Water	04/12/13 09:25	04/12/13 16:34
LL_711_041213	T130856-02	Water	04/12/13 10:20	04/12/13 16:34
LL_715_041213	T130856-03	Water	04/12/13 12:15	04/12/13 16:34
LL_712_041213	T130856-04	Water	04/12/13 14:20	04/12/13 16:34
LL_TB_041213	T130856-05	Water	04/12/13 00:00	04/12/13 16:34

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Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

LL_710_041213
T130856-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	130	50	ug/l	1	3041527	04/15/13	04/16/13	EPA 8015C
Surrogate: 4-Bromofluorobenzene		124 %	65-135		"	"	"	"

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041526	04/15/13	04/16/13	EPA 8260B
Bromochloromethane	ND	1.0	"	"	"	"	"	"
Bromodichloromethane	ND	1.0	"	"	"	"	"	"
Bromoform	ND	1.0	"	"	"	"	"	"
Bromomethane	ND	1.0	"	"	"	"	"	"
n-Butylbenzene	ND	1.0	"	"	"	"	"	"
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"
Chlorobenzene	ND	1.0	"	"	"	"	"	"
Chloroethane	ND	1.0	"	"	"	"	"	"
Chloroform	ND	1.0	"	"	"	"	"	"
Chloromethane	ND	1.0	"	"	"	"	"	"
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"
Dibromochloromethane	ND	1.0	"	"	"	"	"	"
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"
Dibromomethane	ND	1.0	"	"	"	"	"	"
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethane	3.6	1.0	"	"	"	"	"	"
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethene	41	1.0	"	"	"	"	"	"
cis-1,2-Dichloroethene	16	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	1.7	1.0	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Wendy Hsiao

Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

LL_710_041213
T130856-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	3041526	04/15/13	04/16/13	EPA 8260B
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	75	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	89	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"

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Wendy Hsiao

Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

LL_710_041213
T130856-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-butyl alcohol	ND	10	ug/l	1	3041526	04/15/13	04/16/13	EPA 8260B
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		111 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		102 %	81-136		"	"	"	"
Surrogate: Toluene-d8		107 %	88.8-117		"	"	"	"

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Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/19/13 15:40
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LL_711_041213
T130856-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	25000	50	ug/l	1	3041527	04/15/13	04/16/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		85.3 %	65-135		"	"	"	"	

Metals by SM 3500 Series Methods

Ferrous Iron	0.734	0.100	mg/l	1	3041618	04/15/13	04/16/13	EPA6010/S M3500	
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Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041526	04/15/13	04/16/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	27	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	2.6	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	

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Wendy Hsiao

Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

LL_711_041213
T130856-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1-Dichloroethene	ND	1.0	ug/l	1	3041526	04/15/13	04/16/13	EPA 8260B
cis-1,2-Dichloroethene	2.5	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	160	100	"	100	"	"	"	"
p-Isopropyltoluene	4.4	1.0	"	1	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	440	100	"	100	"	"	"	"
n-Propylbenzene	270	100	"	"	"	"	"	"
Styrene	ND	1.0	"	1	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	200	100	"	100	"	"	"	"
1,2,4-Trimethylbenzene	640	100	"	"	"	"	"	"
Vinyl chloride	5.9	1.0	"	1	"	"	"	"
Benzene	2000	50	"	100	"	"	"	"
Toluene	1500	50	"	"	"	"	"	"
Ethylbenzene	450	50	"	"	"	"	"	"
m,p-Xylene	2000	100	"	"	"	"	"	"

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/19/13 15:40
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LL_711_041213
T130856-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

o-Xylene	720	50	ug/l	100	3041526	04/15/13	04/16/13	EPA 8260B	
Tert-amyl methyl ether	ND	2.0	"	1	"	"	"	"	
Tert-butyl alcohol	ND	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene		94.1 %	83.5-119		"	"	"	"	
Surrogate: Dibromofluoromethane		99.8 %	81-136		"	"	"	"	
Surrogate: Toluene-d8		96.0 %	88.8-117		"	"	"	"	

Conventional Chemistry Parameters by APHA/EPA/ASTM Methods

Total Alkalinity	840	20	mg/l	1	3041525	04/15/13	04/15/13	EPA 310.1	
Anions by EPA Method 300.0									
Sulfate as SO4	0.932	0.500	mg/l	1	3041507	04/15/13	04/15/13	EPA 300.0	
Nitrate as NO3	ND	0.500	"	"	"	"	"	"	O-07

RSK-175

Methane	3590	10.0	ug/l	10	3041506	04/15/13	04/15/13	RSK-175	
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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

LL_715_041213
T130856-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	ND	50	ug/l	1	3041527	04/15/13	04/16/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		123 %	65-135		"	"	"	"	

Metals by SM 3500 Series Methods

Ferrous Iron	ND	0.100	mg/l	1	3041618	04/15/13	04/16/13	EPA6010/S M3500	
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Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041526	04/15/13	04/16/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

LL_715_041213
T130856-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1-Dichloroethene	ND	1.0	ug/l	1	3041526	04/15/13	04/16/13	EPA 8260B
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

LL_715_041213
T130856-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

m,p-Xylene	ND	1.0	ug/l	1	3041526	04/15/13	04/16/13	EPA 8260B	
o-Xylene	ND	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	ND	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene		113 %	83.5-119		"	"	"	"	
Surrogate: Dibromofluoromethane		101 %	81-136		"	"	"	"	
Surrogate: Toluene-d8		107 %	88.8-117		"	"	"	"	

Conventional Chemistry Parameters by APHA/EPA/ASTM Methods

Total Alkalinity	540	20	mg/l	1	3041525	04/15/13	04/15/13	EPA 310.1	
Anions by EPA Method 300.0									
Sulfate as SO4	138	2.50	mg/l	5	3041507	04/15/13	04/15/13	EPA 300.0	
Nitrate as NO3	ND	0.500	"	1	"	"	04/15/13	"	O-07
RSK-175									
Methane	48.9	1.00	ug/l	1	3041506	04/15/13	04/15/13	RSK-175	

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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-300 Project Manager: Jeremy Squire	Reported: 04/19/13 15:40
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LL_712_041213
T130856-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	5800	50	ug/l	1	3041527	04/15/13	04/16/13	EPA 8015C
Surrogate: 4-Bromofluorobenzene		91.6 %	65-135		"	"	"	"

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041526	04/15/13	04/16/13	EPA 8260B
Bromochloromethane	ND	1.0	"	"	"	"	"	"
Bromodichloromethane	ND	1.0	"	"	"	"	"	"
Bromoform	ND	1.0	"	"	"	"	"	"
Bromomethane	ND	1.0	"	"	"	"	"	"
n-Butylbenzene	8.8	1.0	"	"	"	"	"	"
sec-Butylbenzene	14	1.0	"	"	"	"	"	"
tert-Butylbenzene	1.9	1.0	"	"	"	"	"	"
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"
Chlorobenzene	ND	1.0	"	"	"	"	"	"
Chloroethane	ND	1.0	"	"	"	"	"	"
Chloroform	ND	1.0	"	"	"	"	"	"
Chloromethane	ND	1.0	"	"	"	"	"	"
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"
Dibromochloromethane	ND	1.0	"	"	"	"	"	"
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"
Dibromomethane	ND	1.0	"	"	"	"	"	"
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"
cis-1,2-Dichloroethene	2.4	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

LL_712_041213
T130856-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,3-Dichloropropane	ND	1.0	ug/l	1	3041526	04/15/13	04/16/13	EPA 8260B	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	52	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	2.7	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	36	1.0	"	"	"	"	"	"	
n-Propylbenzene	51	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	46	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	130	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	540	25	"	50	"	"	"	"	
Toluene	56	0.50	"	1	"	"	"	"	
Ethylbenzene	93	0.50	"	"	"	"	"	"	
m,p-Xylene	390	1.0	"	"	"	"	"	"	
o-Xylene	68	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	ND	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

LL_712_041213
T130856-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Ethyl tert-butyl ether	ND	2.0	ug/l	1	3041526	04/15/13	04/16/13	EPA 8260B	
Methyl tert-butyl ether	180	1.0	"	"	"	"	"	"	E-1
<i>Surrogate: 4-Bromofluorobenzene</i>		101 %	83.5-119		"	"	"	"	
<i>Surrogate: Dibromofluoromethane</i>		98.1 %	81-136		"	"	"	"	
<i>Surrogate: Toluene-d8</i>		104 %	88.8-117		"	"	"	"	

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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

LL_TB_041213
T130856-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041526	04/15/13	04/16/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	

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Wendy Hsiao

Wendy Hsiao For Katherine Shields, Jr. Project Manager



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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

LL_TB_041213
T130856-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

cis-1,3-Dichloropropene	ND	0.50	ug/l	1	3041526	04/15/13	04/16/13	EPA 8260B	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	ND	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	ND	1.0	"	"	"	"	"	"	
n-Propylbenzene	ND	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	ND	0.50	"	"	"	"	"	"	
Toluene	ND	0.50	"	"	"	"	"	"	
Ethylbenzene	ND	0.50	"	"	"	"	"	"	
m,p-Xylene	ND	1.0	"	"	"	"	"	"	
o-Xylene	ND	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	ND	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

LL_TB_041213
T130856-05 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Surrogate: 4-Bromofluorobenzene	107 %	83.5-119	3041526	04/15/13	04/16/13	EPA 8260B
Surrogate: Dibromofluoromethane	97.6 %	81-136	"	"	"	"
Surrogate: Toluene-d8	106 %	88.8-117	"	"	"	"

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

Purgeable Petroleum Hydrocarbons by EPA 8015C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041527 - EPA 5030 GC

Blank (3041527-BLK1)

Prepared: 04/15/13 Analyzed: 04/16/13

C6-C12 (GRO)	ND	50	ug/l							
Surrogate 4-Bromofluorobenzene	120		"	100		120	65-135			

LCS (3041527-BS1)

Prepared: 04/15/13 Analyzed: 04/16/13

C6-C12 (GRO)	4610	50	ug/l	5520		83.5	75-125			
Surrogate 4-Bromofluorobenzene	93.4		"	100		93.4	65-135			

Matrix Spike (3041527-MS1)

Source: T130856-01

Prepared: 04/15/13 Analyzed: 04/16/13

C6-C12 (GRO)	5050	50	ug/l	5520	133	89.0	65-135			
Surrogate 4-Bromofluorobenzene	87.8		"	100		87.8	65-135			

Matrix Spike Dup (3041527-MSD1)

Source: T130856-01

Prepared: 04/15/13 Analyzed: 04/16/13

C6-C12 (GRO)	5230	50	ug/l	5520	133	92.3	65-135	3.53	20	
Surrogate 4-Bromofluorobenzene	101		"	100		101	65-135			

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

Metals by SM 3500 Series Methods - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041618 - EPA 3010A

Blank (3041618-BLK1)

Prepared & Analyzed: 04/16/13

Ferrous Iron	ND	0.100	mg/l
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Duplicate (3041618-DUP1)

Source: T130862-01

Prepared & Analyzed: 04/16/13

Ferrous Iron	0.0260	0.100	mg/l	0.0380	37.5	200
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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041526 - EPA 5030 GCMS

Blank (3041526-BLK1)

Prepared: 04/15/13 Analyzed: 04/16/13

Bromobenzene	ND	1.0	ug/l
Bromochloromethane	ND	1.0	"
Bromodichloromethane	ND	1.0	"
Bromoform	ND	1.0	"
Bromomethane	ND	1.0	"
n-Butylbenzene	ND	1.0	"
sec-Butylbenzene	ND	1.0	"
tert-Butylbenzene	ND	1.0	"
Carbon tetrachloride	ND	0.50	"
Chlorobenzene	ND	1.0	"
Chloroethane	ND	1.0	"
Chloroform	ND	1.0	"
Chloromethane	ND	1.0	"
2-Chlorotoluene	ND	1.0	"
4-Chlorotoluene	ND	1.0	"
Dibromochloromethane	ND	1.0	"
1,2-Dibromo-3-chloropropane	ND	1.0	"
1,2-Dibromoethane (EDB)	ND	1.0	"
Dibromomethane	ND	1.0	"
1,2-Dichlorobenzene	ND	1.0	"
1,3-Dichlorobenzene	ND	1.0	"
1,4-Dichlorobenzene	ND	1.0	"
Dichlorodifluoromethane	ND	0.50	"
1,1-Dichloroethane	ND	1.0	"
1,2-Dichloroethane	ND	0.50	"
1,1-Dichloroethene	ND	1.0	"
cis-1,2-Dichloroethene	ND	1.0	"
trans-1,2-Dichloroethene	ND	1.0	"
1,2-Dichloropropane	ND	1.0	"
1,3-Dichloropropane	ND	1.0	"
2,2-Dichloropropane	ND	1.0	"
1,1-Dichloropropene	ND	1.0	"
cis-1,3-Dichloropropene	ND	0.50	"
trans-1,3-Dichloropropene	ND	0.50	"
Hexachlorobutadiene	ND	1.0	"
Isopropylbenzene	ND	1.0	"

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Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041526 - EPA 5030 GCMS

Blank (3041526-BLK1)

Prepared: 04/15/13 Analyzed: 04/16/13

p-Isopropyltoluene	ND	1.0	ug/l							
Methylene chloride	ND	1.0	"							
Naphthalene	ND	1.0	"							
n-Propylbenzene	ND	1.0	"							
Styrene	ND	1.0	"							
1,1,2,2-Tetrachloroethane	ND	1.0	"							
1,1,1,2-Tetrachloroethane	ND	1.0	"							
Tetrachloroethene	ND	1.0	"							
1,2,3-Trichlorobenzene	ND	1.0	"							
1,2,4-Trichlorobenzene	ND	1.0	"							
1,1,2-Trichloroethane	ND	1.0	"							
1,1,1-Trichloroethane	ND	1.0	"							
Trichloroethene	ND	1.0	"							
Trichlorofluoromethane	ND	1.0	"							
1,2,3-Trichloropropane	ND	1.0	"							
1,3,5-Trimethylbenzene	ND	1.0	"							
1,2,4-Trimethylbenzene	ND	1.0	"							
Vinyl chloride	ND	1.0	"							
Benzene	ND	0.50	"							
Toluene	ND	0.50	"							
Ethylbenzene	ND	0.50	"							
m,p-Xylene	ND	1.0	"							
o-Xylene	ND	0.50	"							
Tert-amyl methyl ether	ND	2.0	"							
Tert-butyl alcohol	ND	10	"							
Di-isopropyl ether	ND	2.0	"							
Ethyl tert-butyl ether	ND	2.0	"							
Methyl tert-butyl ether	ND	1.0	"							
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"							
Surrogate 4-Bromofluorobenzene	8.26		"	8.00		103	83.5-119			
Surrogate Dibromofluoromethane	7.73		"	8.00		96.6	81-136			
Surrogate Toluene-d8	8.51		"	8.00		106	88.8-117			

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Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041526 - EPA 5030 GCMS

LCS (3041526-BS1)

Prepared: 04/15/13 Analyzed: 04/16/13

Chlorobenzene	20.0	1.0	ug/l	20.0		99.8	75-125			
1,1-Dichloroethene	22.2	1.0	"	20.0		111	75-125			
Trichloroethene	20.8	1.0	"	20.0		104	75-125			
Benzene	21.1	0.50	"	20.0		106	75-125			
Toluene	19.7	0.50	"	20.0		98.7	75-125			
Surrogate 4-Bromofluorobenzene	8.47		"	8.00		106	83.5-119			
Surrogate Dibromofluoromethane	8.25		"	8.00		103	81-136			
Surrogate Toluene-d8	8.32		"	8.00		104	88.8-117			

Matrix Spike (3041526-MS1)

Source: T130856-03

Prepared: 04/15/13 Analyzed: 04/16/13

Chlorobenzene	19.6	1.0	ug/l	20.0	ND	98.0	75-125			
1,1-Dichloroethene	22.5	1.0	"	20.0	ND	113	75-125			
Trichloroethene	20.2	1.0	"	20.0	ND	101	75-125			
Benzene	21.5	0.50	"	20.0	ND	108	75-125			
Toluene	19.8	0.50	"	20.0	ND	99.1	75-125			
Surrogate 4-Bromofluorobenzene	8.37		"	8.00		105	83.5-119			
Surrogate Dibromofluoromethane	8.67		"	8.00		108	81-136			
Surrogate Toluene-d8	8.29		"	8.00		104	88.8-117			

Matrix Spike Dup (3041526-MSD1)

Source: T130856-03

Prepared: 04/15/13 Analyzed: 04/16/13

Chlorobenzene	19.4	1.0	ug/l	20.0	ND	97.2	75-125	0.717	20	
1,1-Dichloroethene	22.2	1.0	"	20.0	ND	111	75-125	1.38	20	
Trichloroethene	20.3	1.0	"	20.0	ND	102	75-125	0.444	20	
Benzene	21.4	0.50	"	20.0	ND	107	75-125	0.233	20	
Toluene	19.9	0.50	"	20.0	ND	99.6	75-125	0.553	20	
Surrogate 4-Bromofluorobenzene	8.55		"	8.00		107	83.5-119			
Surrogate Dibromofluoromethane	8.55		"	8.00		107	81-136			
Surrogate Toluene-d8	8.40		"	8.00		105	88.8-117			

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Conventional Chemistry Parameters by APHA/EPA/ASTM Methods - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041525 - General Preparation

Duplicate (3041525-DUP1) **Source: T130856-03** Prepared & Analyzed: 04/15/13

Total Alkalinity	555	20	mg/l		545			1.82	25	
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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

Anions by EPA Method 300.0 - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041507 - General Preparation

Blank (3041507-BLK1)

Prepared & Analyzed: 04/15/13

Sulfate as SO ₄	ND	0.500	mg/l
Nitrate as NO ₃	ND	0.500	"

LCS (3041507-BS1)

Prepared & Analyzed: 04/15/13

Sulfate as SO ₄	10.5	0.500	mg/l	10.0	105	75-125
Nitrate as NO ₃	0.574	0.500	"	0.500	115	75-125

Matrix Spike (3041507-MS1)

Source: T130856-02

Prepared & Analyzed: 04/15/13

Sulfate as SO ₄	10.5	0.500	mg/l	10.0	0.932	96.1	75-125
Nitrate as NO ₃	0.585	0.500	"	0.500	ND	117	75-125

Matrix Spike Dup (3041507-MSD1)

Source: T130856-02

Prepared & Analyzed: 04/15/13

Sulfate as SO ₄	10.2	0.500	mg/l	10.0	0.932	93.0	75-125	2.96	20
Nitrate as NO ₃	0.566	0.500	"	0.500	ND	113	75-125	3.30	20

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Project: Cenco
Project Number: 1003-001-300
Project Manager: Jeremy Squire

Reported:
04/19/13 15:40

RSK-175 - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041506 - EPA 3810m Headspace

Blank (3041506-BLK1)

Prepared & Analyzed: 04/15/13

Methane	ND	1.00	ug/l
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Duplicate (3041506-DUP1)

Source: T130856-02

Prepared & Analyzed: 04/15/13

Methane	3380	10.0	ug/l	3590	6.06	20
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Notes and Definitions

O-07 The sample was analyzed outside the EPA recommended holding time of 48 hours.

E-1 The final dilution was lower than the original data or previous dilutions. The highest recovered concentration was reported even though it was above calibration range.

DET Analyte DETECTED

ND Analyte NOT DETECTED at or above the reporting limit

NR Not Reported

dry Sample results reported on a dry weight basis

RPD Relative Percent Difference

SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Wendy Hsiao For Katherine Shields, Jr. Project Manager

SunStar Laboratories, Inc.
25712 Commercentre Dr
Lake Forest, CA 92630
949-297-5020

Chain of Custody Record

Client: MUREX ENVIRONMENTAL INC.
Address: 2640 Walnut Ave, Unit F
Phone: (714) 508-0800 Fax: (714) 508-0880
Project Manager: Jeremy Squire (714) 604-5836

Date: 04/12/2013

Page: 1 OF 1

Project Name: CENCO

Collector: ~~Franco Sosa~~ AW

Client Project #: 1003-001-300

Batch #: 7130856

EDF #: _____

Sample ID	Date Sampled	Time	Sample Type	TPHg (8015 M)	VOCs (8260 B)	METHANE (8015)	ALKALINITY	3001C (NITRATE, SULFATE)	FERROUS IRON									Total # of containers	Comments/Preservative	Laboratory ID #
LL-710-041213	04/12/13	0925	GW	X	X													6		01
LL-711-041213	↓	1020	↓	X	X	X	X	X	X									12		02
LL-715-041213	↓	1215	↓	X	X	X	X	X	X									12		03
LL-712-041213	↓	1420	↓	X	X													6		04
LL-TB-041213	—	—	WATER															2		05
Relinquished by: (signature) <i>T.A.L.</i>	Date / Time 04/12/13 16:34	Received by: (Sign / Date / Time) <i>[Signature]</i> 4-12-13 16:34	Total # of containers	38	Notes															
Relinquished by: (signature)	Date / Time	Received by: (Sign / Date / Time)	Chain of Custody seals																	
Relinquished by: (signature)	Date / Time	Received by: (Sign / Date / Time)	Seals intact? Y/N/NA																	
Relinquished by: (signature)	Date / Time	Received by: (Sign / Date / Time)	Received good condition/cold	4.6																
Relinquished by: (signature)	Date / Time	Received by: (Sign / Date / Time)	Turn around time:	Standard																

Sample disposal Instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____

SAMPLE RECEIVING REVIEW SHEET

BATCH # 7130856

Client Name: MUREX

Project: CENCO

Received by: SUNNY

Date/Time Received: 4.12.13 / 16:34

Delivered by: ☐ Client ☒ SunStar Courier ☐ GSO ☐ FedEx ☐ Other _____

Total number of coolers received 0 Temp criteria = 6°C > 0°C (no frozen containers)

Temperature: cooler #1 4.8 °C +/- the CF (- 0.2°C) = 4.6 °C corrected temperature

cooler #2 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

cooler #3 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

Samples outside temp. but received on ice, w/in 6 hours of final sampling. ☒ Yes ☐ No* ☐ N/A

Custody Seals Intact on Cooler/Sample ☐ Yes ☐ No* ☒ N/A

Sample Containers Intact ☒ Yes ☐ No*

Sample labels match COC ID's ☒ Yes ☐ No*

Total number of containers received match COC ☒ Yes ☐ No*

Proper containers received for analyses requested on COC ☒ Yes ☐ No*

Proper preservative indicated on COC/containers for analyses requested ☒ Yes ☐ No* ☐ N/A

Complete shipment received in good condition with correct temperatures, containers, labels, volumes preservatives and within method specified holding times. ☒ Yes ☐ No*

* Complete Non-Conformance Receiving Sheet if checked

Cooler/Sample Review - Initials and date SI 4.12.13

Comments:



25712 Commercentre Drive
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22 April 2013

Jeremy Squire
Murex
15375 Barranca Parkway, Suite K-101
Irvine, CA 92861
RE: Cenco

Enclosed are the results of analyses for samples received by the laboratory on 04/15/13 15:15. If you have any questions concerning this report, please feel free to contact me.

Sincerely,

Katherine Shields
Jr. Project Manager



25712 Commercentre Drive
Lake Forest, California 92630
949.297.5020 Phone
949.297.5027 Fax

Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
LL_706_041513_01	T130862-01	Water	04/15/13 09:20	04/15/13 15:15
LL_706_041513_02	T130862-02	Water	04/15/13 09:30	04/15/13 15:15
LL_704_041513	T130862-03	Water	04/15/13 11:30	04/15/13 15:15
LL_TB_041513	T130862-04	Water	04/15/13 11:30	04/15/13 15:15

SunStar Laboratories, Inc.

Katherine Shields

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Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

LL_706_041513_01
T130862-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	260	50	ug/l	1	3041617	04/16/13	04/17/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		109 %	65-135		"	"	"	"	

Metals by SM 3500 Series Methods

Ferrous Iron	ND	0.100	mg/l	1	3041618	04/16/13	04/16/13	EPA6010/S M3500	
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Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041616	04/16/13	04/17/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

LL_706_041513_01
T130862-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1-Dichloroethene	ND	1.0	ug/l	1	3041616	04/16/13	04/17/13	EPA 8260B
cis-1,2-Dichloroethene	4.0	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	1.0	1.0	"	"	"	"	"	"
Benzene	5.9	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-200 Project Manager: Jeremy Squire	Reported: 04/22/13 15:43
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LL_706_041513_01
T130862-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

m,p-Xylene	ND	1.0	ug/l	1	3041616	04/16/13	04/17/13	EPA 8260B
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"
Tert-butyl alcohol	54	10	"	"	"	"	"	"
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	2.8	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		108 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		99.9 %	81-136		"	"	"	"
Surrogate: Toluene-d8		105 %	88.8-117		"	"	"	"

Conventional Chemistry Parameters by APHA/EPA/ASTM Methods

Total Alkalinity	570	20	mg/l	1	3041614	04/16/13	04/16/13	EPA 310.1
Anions by EPA Method 300.0								
Sulfate as SO4	77.9	1.00	mg/l	2	3041611	04/16/13	04/16/13	EPA 300.0
Nitrate as NO3	ND	0.500	"	1	"	"	04/16/13	"
RSK-175								
Methane	461	10.0	ug/l	10	3041619	04/16/13	04/18/13	RSK-175

SunStar Laboratories, Inc.

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15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

LL_706_041513_02
T130862-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	250	50	ug/l	1	3041617	04/16/13	04/17/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		105 %	65-135		"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041616	04/16/13	04/17/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	4.9	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

LL_706_041513_02
T130862-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,3-Dichloropropane	ND	1.0	ug/l	1	3041616	04/16/13	04/17/13	EPA 8260B
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	ND	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	1.5	1.0	"	"	"	"	"	"
Benzene	5.1	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"
Tert-butyl alcohol	61	10	"	"	"	"	"	"

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager

Page 6 of 22



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

LL_706_041513_02
T130862-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Di-isopropyl ether	ND	2.0	ug/l	1	3041616	04/16/13	04/17/13	EPA 8260B
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	3.2	1.0	"	"	"	"	"	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene		109 %	83.5-119		"	"	"	"
Surrogate: Dibromofluoromethane		102 %	81-136		"	"	"	"
Surrogate: Toluene-d8		104 %	88.8-117		"	"	"	"

SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

LL_704_041513
T130862-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Purgeable Petroleum Hydrocarbons by EPA 8015C

C6-C12 (GRO)	3900	50	ug/l	1	3041617	04/16/13	04/17/13	EPA 8015C	
Surrogate: 4-Bromofluorobenzene		90.4 %	65-135		"	"	"	"	

Metals by SM 3500 Series Methods

Ferrous Iron	0.201	0.100	mg/l	1	3041618	04/16/13	04/16/13	EPA6010/S M3500	
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Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041616	04/16/13	04/17/13	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	7.9	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	2.4	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	3.2	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	6.1	0.50	"	"	"	"	"	"	

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Katherine Shields

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Murex
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Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

LL_704_041513
T130862-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1-Dichloroethene	ND	1.0	ug/l	1	3041616	04/16/13	04/17/13	EPA 8260B	
cis-1,2-Dichloroethene	3.0	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	75	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	12	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	57	1.0	"	"	"	"	"	"	
n-Propylbenzene	83	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	170	25	"	25	"	"	"	"	
1,2,4-Trimethylbenzene	530	25	"	"	"	"	"	"	
Vinyl chloride	1.1	1.0	"	1	"	"	"	"	
Benzene	420	0.50	"	"	"	"	"	"	E-1
Toluene	29	0.50	"	"	"	"	"	"	
Ethylbenzene	200	12	"	25	"	"	"	"	
m,p-Xylene	300	25	"	"	"	"	"	"	

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Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

LL_704_041513
T130862-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

o-Xylene	10	0.50	ug/l	1	3041616	04/16/13	04/17/13	EPA 8260B	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	ND	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	97	1.0	"	"	"	"	"	"	
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene		97.2 %	83.5-119		"	"	"	"	
Surrogate: Dibromofluoromethane		104 %	81-136		"	"	"	"	
Surrogate: Toluene-d8		104 %	88.8-117		"	"	"	"	

Conventional Chemistry Parameters by APHA/EPA/ASTM Methods

Total Alkalinity	820	20	mg/l	1	3041614	04/16/13	04/16/13	EPA 310.1	
Anions by EPA Method 300.0									
Sulfate as SO4	16.4	0.500	mg/l	1	3041611	04/16/13	04/16/13	EPA 300.0	
Nitrate as NO3	ND	0.500	"	"	"	"	"	"	
RSK-175									
Methane	1970	25.0	ug/l	25	3041619	04/16/13	04/18/13	RSK-175	

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Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

LL_TB_041513
T130862-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	3041616	04/16/13	04/17/13	EPA 8260B
Bromochloromethane	ND	1.0	"	"	"	"	"	"
Bromodichloromethane	ND	1.0	"	"	"	"	"	"
Bromoform	ND	1.0	"	"	"	"	"	"
Bromomethane	ND	1.0	"	"	"	"	"	"
n-Butylbenzene	ND	1.0	"	"	"	"	"	"
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"
Chlorobenzene	ND	1.0	"	"	"	"	"	"
Chloroethane	ND	1.0	"	"	"	"	"	"
Chloroform	ND	1.0	"	"	"	"	"	"
Chloromethane	ND	1.0	"	"	"	"	"	"
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"
Dibromochloromethane	ND	1.0	"	"	"	"	"	"
1,2-Dibromo-3-chloropropane	ND	1.0	"	"	"	"	"	"
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"
Dibromomethane	ND	1.0	"	"	"	"	"	"
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

LL_TB_041513
T130862-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

cis-1,3-Dichloropropene	ND	0.50	ug/l	1	3041616	04/16/13	04/17/13	EPA 8260B
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"
Isopropylbenzene	ND	1.0	"	"	"	"	"	"
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"
Methylene chloride	ND	1.0	"	"	"	"	"	"
Naphthalene	1.2	1.0	"	"	"	"	"	"
n-Propylbenzene	ND	1.0	"	"	"	"	"	"
Styrene	ND	1.0	"	"	"	"	"	"
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"
Tetrachloroethene	ND	1.0	"	"	"	"	"	"
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"
Trichloroethene	ND	1.0	"	"	"	"	"	"
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"
Vinyl chloride	ND	1.0	"	"	"	"	"	"
Benzene	ND	0.50	"	"	"	"	"	"
Toluene	ND	0.50	"	"	"	"	"	"
Ethylbenzene	ND	0.50	"	"	"	"	"	"
m,p-Xylene	ND	1.0	"	"	"	"	"	"
o-Xylene	ND	0.50	"	"	"	"	"	"
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"
Tert-butyl alcohol	ND	10	"	"	"	"	"	"
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"

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Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

LL_TB_041513
T130862-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	ug/l	1	3041616	04/16/13	04/17/13	EPA 8260B	
Surrogate: 4-Bromofluorobenzene	107 %	83.5-119			"	"	"	"	
Surrogate: Dibromofluoromethane	105 %	81-136			"	"	"	"	
Surrogate: Toluene-d8	105 %	88.8-117			"	"	"	"	

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Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

Purgeable Petroleum Hydrocarbons by EPA 8015C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041617 - EPA 5030 GC

Blank (3041617-BLK1)

Prepared: 04/16/13 Analyzed: 04/17/13

C6-C12 (GRO)	ND	50	ug/l							
Surrogate 4-Bromofluorobenzene	117		"	100		117	65-135			

LCS (3041617-BS1)

Prepared: 04/16/13 Analyzed: 04/17/13

C6-C12 (GRO)	5470	50	ug/l	5520		99.1	75-125			
Surrogate 4-Bromofluorobenzene	70.2		"	100		70.2	65-135			

Matrix Spike (3041617-MS1)

Source: T130862-01

Prepared: 04/16/13 Analyzed: 04/17/13

C6-C12 (GRO)	5770	50	ug/l	5520	265	99.8	65-135			
Surrogate 4-Bromofluorobenzene	74.8		"	100		74.8	65-135			

Matrix Spike Dup (3041617-MSD1)

Source: T130862-01

Prepared: 04/16/13 Analyzed: 04/17/13

C6-C12 (GRO)	5210	50	ug/l	5520	265	89.7	65-135	10.2	20	
Surrogate 4-Bromofluorobenzene	71.7		"	100		71.7	65-135			

SunStar Laboratories, Inc.

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Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

Metals by SM 3500 Series Methods - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041618 - EPA 3010A

Blank (3041618-BLK1)

Prepared & Analyzed: 04/16/13

Ferrous Iron	ND	0.100	mg/l
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Duplicate (3041618-DUP1)

Source: T130862-01

Prepared & Analyzed: 04/16/13

Ferrous Iron	0.0260	0.100	mg/l	0.0380	37.5	200
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Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041616 - EPA 5030 GCMS

Blank (3041616-BLK1)

Prepared & Analyzed: 04/16/13

Bromobenzene	ND	1.0	ug/l
Bromochloromethane	ND	1.0	"
Bromodichloromethane	ND	1.0	"
Bromoform	ND	1.0	"
Bromomethane	ND	1.0	"
n-Butylbenzene	ND	1.0	"
sec-Butylbenzene	ND	1.0	"
tert-Butylbenzene	ND	1.0	"
Carbon tetrachloride	ND	0.50	"
Chlorobenzene	ND	1.0	"
Chloroethane	ND	1.0	"
Chloroform	ND	1.0	"
Chloromethane	ND	1.0	"
2-Chlorotoluene	ND	1.0	"
4-Chlorotoluene	ND	1.0	"
Dibromochloromethane	ND	1.0	"
1,2-Dibromo-3-chloropropane	ND	1.0	"
1,2-Dibromoethane (EDB)	ND	1.0	"
Dibromomethane	ND	1.0	"
1,2-Dichlorobenzene	ND	1.0	"
1,3-Dichlorobenzene	ND	1.0	"
1,4-Dichlorobenzene	ND	1.0	"
Dichlorodifluoromethane	ND	0.50	"
1,1-Dichloroethane	ND	1.0	"
1,2-Dichloroethane	ND	0.50	"
1,1-Dichloroethene	ND	1.0	"
cis-1,2-Dichloroethene	ND	1.0	"
trans-1,2-Dichloroethene	ND	1.0	"
1,2-Dichloropropane	ND	1.0	"
1,3-Dichloropropane	ND	1.0	"
2,2-Dichloropropane	ND	1.0	"
1,1-Dichloropropene	ND	1.0	"
cis-1,3-Dichloropropene	ND	0.50	"
trans-1,3-Dichloropropene	ND	0.50	"
Hexachlorobutadiene	ND	1.0	"
Isopropylbenzene	ND	1.0	"

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Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041616 - EPA 5030 GCMS

Blank (3041616-BLK1)

Prepared & Analyzed: 04/16/13

p-Isopropyltoluene	ND	1.0	ug/l
Methylene chloride	ND	1.0	"
Naphthalene	ND	1.0	"
n-Propylbenzene	ND	1.0	"
Styrene	ND	1.0	"
1,1,2,2-Tetrachloroethane	ND	1.0	"
1,1,1,2-Tetrachloroethane	ND	1.0	"
Tetrachloroethene	ND	1.0	"
1,2,3-Trichlorobenzene	ND	1.0	"
1,2,4-Trichlorobenzene	ND	1.0	"
1,1,2-Trichloroethane	ND	1.0	"
1,1,1-Trichloroethane	ND	1.0	"
Trichloroethene	ND	1.0	"
Trichlorofluoromethane	ND	1.0	"
1,2,3-Trichloropropane	ND	1.0	"
1,3,5-Trimethylbenzene	ND	1.0	"
1,2,4-Trimethylbenzene	ND	1.0	"
Vinyl chloride	ND	1.0	"
Benzene	ND	0.50	"
Toluene	ND	0.50	"
Ethylbenzene	ND	0.50	"
m,p-Xylene	ND	1.0	"
o-Xylene	ND	0.50	"
Tert-amyl methyl ether	ND	2.0	"
Tert-butyl alcohol	ND	10	"
Di-isopropyl ether	ND	2.0	"
Ethyl tert-butyl ether	ND	2.0	"
Methyl tert-butyl ether	ND	1.0	"
1,1,2-trichloro-1,2,2-trifluoroethane (CFC 113)	ND	5.0	"
Surrogate 4-Bromofluorobenzene	8.73		"
Surrogate Dibromofluoromethane	8.25		"
Surrogate Toluene-d8	8.41		"

8.00 109 83.5-119
8.00 103 81-136
8.00 105 88.8-117

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Murex	Project: Cenco	Reported:
15375 Barranca Parkway, Suite K-101	Project Number: 1003-001-200	04/22/13 15:43
Irvine CA, 92861	Project Manager: Jeremy Squire	

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041616 - EPA 5030 GCMS

LCS (3041616-BS1)

Prepared: 04/16/13 Analyzed: 04/17/13

Chlorobenzene	20.1	1.0	ug/l	20.0		100	75-125			
1,1-Dichloroethene	25.2	1.0	"	20.0		126	75-125			QM-13
Trichloroethene	23.2	1.0	"	20.0		116	75-125			
Benzene	23.8	0.50	"	20.0		119	75-125			
Toluene	19.0	0.50	"	20.0		95.2	75-125			
Surrogate 4-Bromofluorobenzene	8.44		"	8.00		106	83.5-119			
Surrogate Dibromofluoromethane	9.13		"	8.00		114	81-136			
Surrogate Toluene-d8	8.39		"	8.00		105	88.8-117			

Matrix Spike (3041616-MS1)

Source: T130862-01

Prepared: 04/16/13 Analyzed: 04/17/13

Chlorobenzene	19.2	1.0	ug/l	20.0	ND	96.0	75-125			
1,1-Dichloroethene	24.1	1.0	"	20.0	ND	121	75-125			
Trichloroethene	21.3	1.0	"	20.0	ND	106	75-125			
Benzene	29.1	0.50	"	20.0	5.94	116	75-125			
Toluene	20.2	0.50	"	20.0	0.350	99.5	75-125			
Surrogate 4-Bromofluorobenzene	8.31		"	8.00		104	83.5-119			
Surrogate Dibromofluoromethane	9.22		"	8.00		115	81-136			
Surrogate Toluene-d8	8.30		"	8.00		104	88.8-117			

Matrix Spike Dup (3041616-MSD1)

Source: T130862-01

Prepared: 04/16/13 Analyzed: 04/17/13

Chlorobenzene	21.2	1.0	ug/l	20.0	ND	106	75-125	10.0	20	
1,1-Dichloroethene	26.1	1.0	"	20.0	ND	130	75-125	7.76	20	QM-13
Trichloroethene	21.8	1.0	"	20.0	ND	109	75-125	2.23	20	
Benzene	30.8	0.50	"	20.0	5.94	124	75-125	5.51	20	
Toluene	21.0	0.50	"	20.0	0.350	104	75-125	3.87	20	
Surrogate 4-Bromofluorobenzene	8.59		"	8.00		107	83.5-119			
Surrogate Dibromofluoromethane	9.49		"	8.00		119	81-136			
Surrogate Toluene-d8	8.13		"	8.00		102	88.8-117			

SunStar Laboratories, Inc.

The results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex 15375 Barranca Parkway, Suite K-101 Irvine CA, 92861	Project: Cenco Project Number: 1003-001-200 Project Manager: Jeremy Squire	Reported: 04/22/13 15:43
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Conventional Chemistry Parameters by APHA/EPA/ASTM Methods - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041614 - General Preparation

Duplicate (3041614-DUP1) **Source: T130862-01** Prepared & Analyzed: 04/16/13

Total Alkalinity	595	20	mg/l	570	4.29	25
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SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager



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Murex	Project: Cenco	Reported:
15375 Barranca Parkway, Suite K-101	Project Number: 1003-001-200	04/22/13 15:43
Irvine CA, 92861	Project Manager: Jeremy Squire	

Anions by EPA Method 300.0 - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041611 - General Preparation

Blank (3041611-BLK1)				Prepared & Analyzed: 04/16/13						
Sulfate as SO4	ND	0.500	mg/l							
Nitrate as NO3	ND	0.500	"							
LCS (3041611-BS1)				Prepared & Analyzed: 04/16/13						
Sulfate as SO4	9.88	0.500	mg/l	10.0		98.8	75-125			
Nitrate as NO3	0.531	0.500	"	0.500		106	75-125			
Matrix Spike (3041611-MS1)				Source: T130862-01		Prepared & Analyzed: 04/16/13				
Sulfate as SO4	84.2	0.500	mg/l	10.0	77.9	62.9	75-125			QM-02
Nitrate as NO3	0.601	0.500	"	0.500	ND	120	75-125			
Matrix Spike Dup (3041611-MSD1)				Source: T130862-01		Prepared & Analyzed: 04/16/13				
Sulfate as SO4	84.8	0.500	mg/l	10.0	77.9	69.0	75-125	0.726	20	QM-02
Nitrate as NO3	0.635	0.500	"	0.500	ND	127	75-125	5.50	20	QM-05

SunStar Laboratories, Inc.

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Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

RSK-175 - Quality Control
SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 3041619 - EPA 3810m Headspace

Blank (3041619-BLK1)

Prepared: 04/16/13 Analyzed: 04/18/13

Methane	ND	1.00	ug/l
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Duplicate (3041619-DUP1)

Source: T130862-01

Prepared: 04/16/13 Analyzed: 04/18/13

Methane	486	10.0	ug/l	461	5.11	20
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SunStar Laboratories, Inc.

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Katherine Shields

Katherine Shields, Jr. Project Manager

Murex
15375 Barranca Parkway, Suite K-101
Irvine CA, 92861

Project: Cenco
Project Number: 1003-001-200
Project Manager: Jeremy Squire

Reported:
04/22/13 15:43

Notes and Definitions

- QM-13 Spike recovery for this analyte was bias high in the LCS and/or MS/MSD. Instrument blank, method blank and all samples were ND. No negative impact on data is expected.
- QM-05 The spike recovery was outside acceptance limits for the MS and/or MSD due to possible matrix interference. The LCS was within acceptance criteria. The data is acceptable as no negative impact on data is expected.
- QM-02 The RPD and/or percent recovery for this QC spike sample cannot be accurately calculated due to the high concentration of analyte inherent in the sample.
- E-1 The final dilution was lower than the original data or previous dilutions. The highest recovered concentration was reported even though it was above calibration range.
- DET Analyte DETECTED
- ND Analyte NOT DETECTED at or above the reporting limit
- NR Not Reported
- dry Sample results reported on a dry weight basis
- RPD Relative Percent Difference

Chain of Custody Record

Date: 04/15/2013

Page: 1 OF 1

Project Name: CENCO

Collector: ~~Frano Sosic~~ *AW*

Client Project #: 1003-001-300

Batch #: T130862

EDF #:

Sample ID	Date Sampled	Time	Sample Type	TPHg (8015 M)	VOCs (8260 B)	METHANE (8015)	ALKALINITY	300 IC (NITRATE, SULFATE)	FERROUS (PbZn)								Total # of containers	Comments/Preservative	Laboratory ID #
LL-706-041513-01	04/15/13	0920	GW	X	X	X	X	X	X								12		01
LL-706-041513-02	↓	0930	↓	X	X												6		02
LL-704-041513	↓	1130	↓	X	X	X	X	X	X								12		03
LL-TB-041513	-	-	WATER	X													2		04
Relinquished by: (signature) T.A.W.	Date / Time 04/15/13	1515	Received by: (Sign / Date / Time) <i>[Signature]</i> 4/15/13 1515	Total # of containers			Chain of Custody seals			Seals intact? Y/N/NA			Received good condition/cold			Turn around time:			Notes
Relinquished by: (signature)	Date / Time		Received by: (Sign / Date / Time)	32			NA			NA			5.6			Standard			
Relinquished by: (signature)	Date / Time		Received by: (Sign / Date / Time)																

Sample disposal Instructions: Disposal @ \$2.00 each _____ Return to client _____ Pickup _____

SAMPLE RECEIVING REVIEW SHEET

BATCH # T130862

Client Name: MUREX

Project: CENCO

Received by: PATRICK

Date/Time Received: 4/15/13 1515

Delivered by: ☐ Client ☒ SunStar Courier ☐ GSO ☐ FedEx ☐ Other _____

Total number of coolers received 0

Temp criteria = 6°C > 0°C (no frozen containers)

Temperature: cooler #1 5.8 °C +/- the CF (- 0.2°C) = 5.6 °C corrected temperature

cooler #2 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

cooler #3 _____ °C +/- the CF (- 0.2°C) = _____ °C corrected temperature

Samples outside temp. but received on ice, w/in 6 hours of final sampling. ☒ Yes ☐ No* ☐ N/A

Custody Seals Intact on Cooler/Sample ☐ Yes ☐ No* ☒ N/A

Sample Containers Intact ☒ Yes ☐ No*

Sample labels match COC ID's ☒ Yes ☐ No*

Total number of containers received match COC ☒ Yes ☐ No*

Proper containers received for analyses requested on COC ☒ Yes ☐ No*

Proper preservative indicated on COC/containers for analyses requested ☒ Yes ☐ No* ☐ N/A

Complete shipment received in good condition with correct temperatures, containers, labels, volumes preservatives and within method specified holding times. ☒ Yes ☐ No*

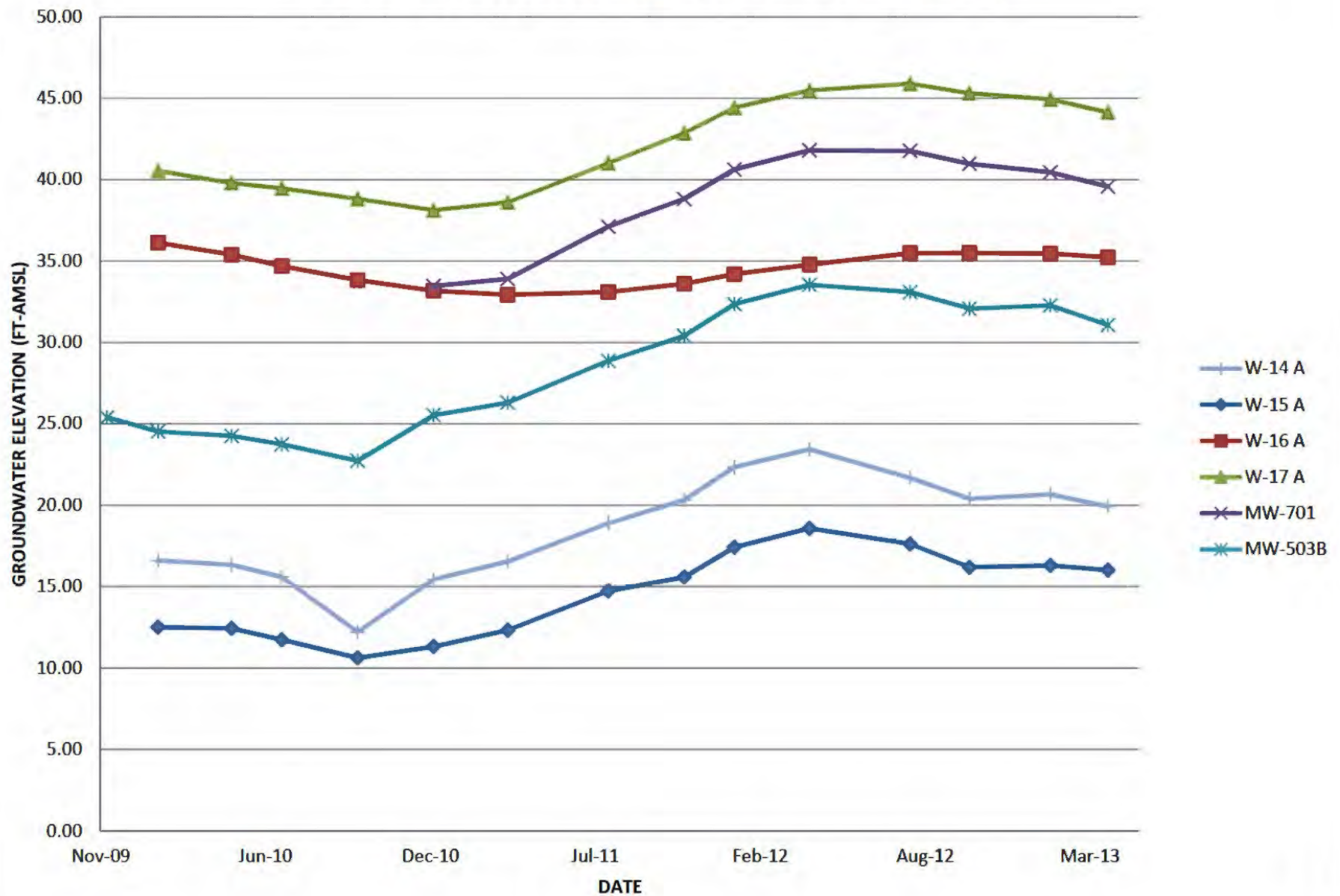
* Complete Non-Conformance Receiving Sheet if checked

Cooler/Sample Review - Initials and date BC 4/15/13

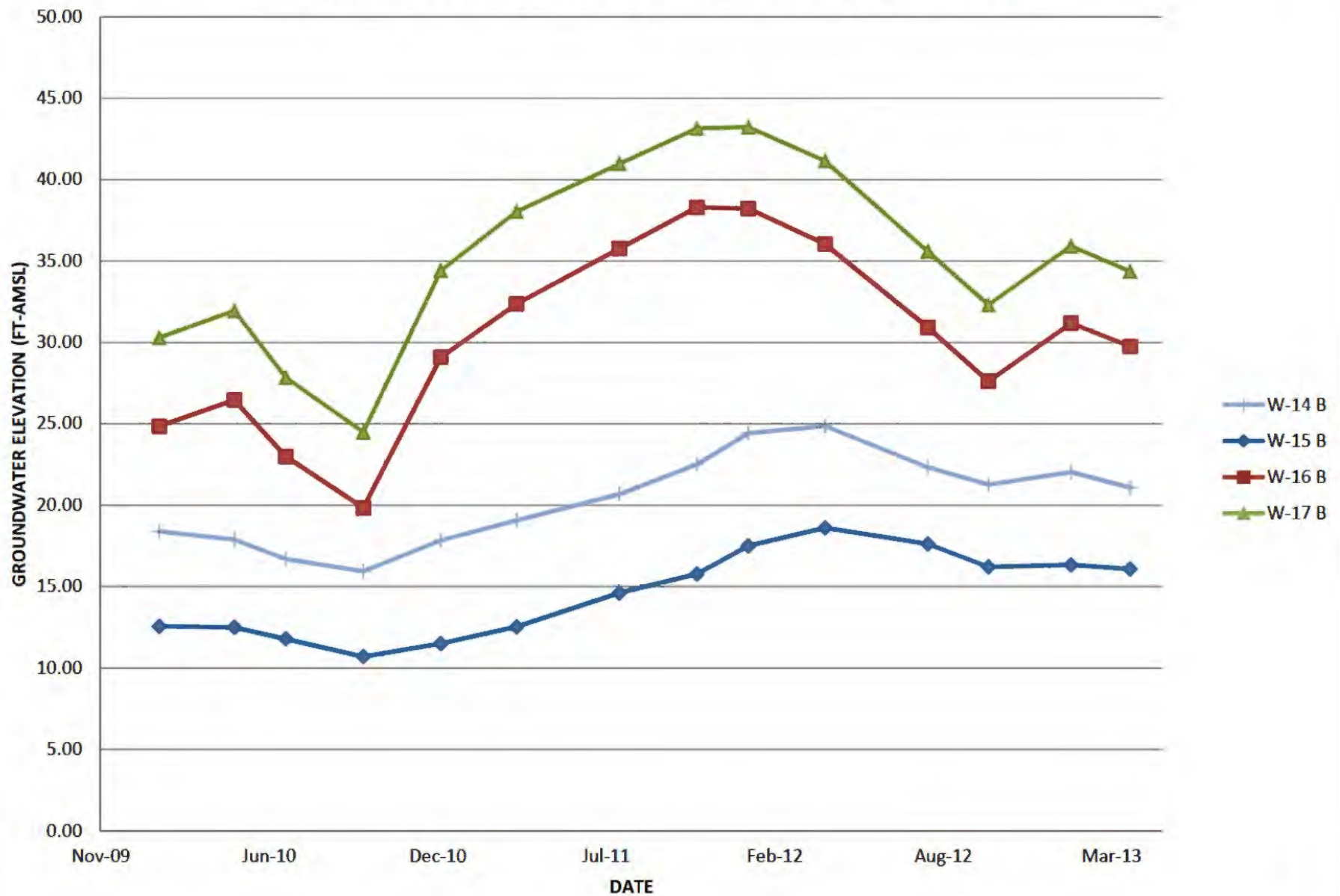
Comments:

Appendix C

A - SCREENED WELLS (APPROX 60-120 FT-BGS)



B - SCREENED WELLS (APPROX 145-170 FT-BGS)



C - SCREENED WELLS (APPROX 185-200 FT-BGS)

